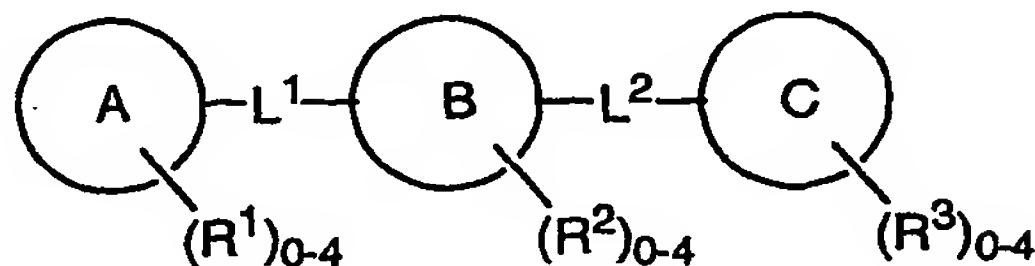


*What is claimed is:*

1. A compound for modulating c-Kit activity according to Formula II,



II

or a pharmaceutically acceptable salt, hydrate, or prodrug thereof, wherein,

ring A is a five- to fourteen-membered heteroaryl;

each R<sup>1</sup> is independently selected from -H, halogen, trihalomethyl, -CN, -NO<sub>2</sub>, -OR<sup>4</sup>, -N(R<sup>4</sup>)R<sup>4</sup>, -S(O)<sub>0-2</sub>R<sup>4</sup>, -SO<sub>2</sub>N(R<sup>4</sup>)R<sup>4</sup>, -CO<sub>2</sub>R<sup>4</sup>, -C(=O)N(R<sup>4</sup>)R<sup>4</sup>, -C(=NR<sup>5</sup>)N(R<sup>4</sup>)R<sup>4</sup>, -C(=NR<sup>5</sup>)R<sup>4</sup>, -N(R<sup>4</sup>)SO<sub>2</sub>R<sup>4</sup>, -N(R<sup>4</sup>)C(O)R<sup>4</sup>, -NCO<sub>2</sub>R<sup>4</sup>, -C(=O)R<sup>4</sup>, optionally substituted alkoxy, optionally substituted C<sub>1-6</sub>alkyl, optionally substituted aryl, optionally substituted aryl C<sub>1-6</sub>alkyl, optionally substituted heterocyclyl, and optionally substituted heterocyclyl C<sub>1-6</sub>alkyl;

two adjacent of R<sup>1</sup>, together with the annular atoms to which they are attached, can form a five- to six-membered ring containing up to two heteroatoms and optionally substituted with up to three of R<sup>10</sup>;

L<sup>1</sup> is selected from a single bond, an optionally substituted C<sub>1-2</sub>alkylene, -O-, -CH<sub>2</sub>O-, -N(R<sup>7</sup>)-, -C(=O)N(R<sup>7</sup>)-, -SO<sub>2</sub>N(R<sup>7</sup>)-, -CH<sub>2</sub>N(R<sup>7</sup>)-, and -S(O)<sub>0-2</sub>-;

ring B is a five- to ten-membered aryl or a five- to ten-membered heterocyclyl;

each R<sup>2</sup> is independently selected from -H, halogen, trihalomethyl, -CN, -NO<sub>2</sub>, -OR<sup>4</sup>, -N(R<sup>4</sup>)R<sup>4</sup>, -S(O)<sub>0-2</sub>R<sup>4</sup>, -SO<sub>2</sub>N(R<sup>4</sup>)R<sup>4</sup>, -CO<sub>2</sub>R<sup>4</sup>, -C(=O)N(R<sup>4</sup>)R<sup>4</sup>, -C(=NR<sup>5</sup>)N(R<sup>4</sup>)R<sup>4</sup>, -C(=NR<sup>5</sup>)R<sup>4</sup>, -N(R<sup>4</sup>)SO<sub>2</sub>R<sup>4</sup>, -N(R<sup>4</sup>)C(O)R<sup>4</sup>, -NCO<sub>2</sub>R<sup>4</sup>, -C(=O)R<sup>4</sup>, optionally substituted alkoxy, optionally substituted C<sub>1-6</sub>alkyl, optionally substituted aryl, optionally substituted aryl C<sub>1-6</sub>alkyl, optionally substituted heterocyclyl, and optionally substituted heterocyclyl C<sub>1-6</sub>alkyl;

two adjacent of R<sup>2</sup>, together with the annular atoms to which they are attached, can form a five- to six-membered ring containing up to two heteroatoms and optionally substituted with up to three of R<sup>15</sup>;

$L^2$  is selected from  $C_4$ alkylene,  $C_4$ alkylidene,  $C_4$ alkylidyne,  $-X(CH_2)_2O-$ ,  $-X(CH_2)_2N(R^7)-$ ,  $-XCH_2SO_2N(R^7)-$ ,  $-XN(R^7)C(=O)N(R^7)-$ ,  $-XCH_2C(=O)N(R^7)-$ ,  $-(CH_2)_3X-$ ,  $-XN(R^7)SO_2N(R^7)-$ ,  $-XCH_2N(R^7)SO_2-$ ,  $-CH_2X(CH_2)_2-$ ,  $-CH=CHC(=O)N(R^7)-$ ,  $-CH=CHSO_2N(R^7)-$ ,  $-XCH_2N(R^7)C(=O)-$ ,  $-M-M-$ ,  $-CH_2N(R^7)C(=O)O-$ , and  $-CH_2OC(=O)N(R^7)-$ ; wherein X is selected from  $-CH_2-$ ,  $-O-$ ,  $-N(R^7)-$ ,  $-C(=O)-$ , and  $-S(O)_{0-2}-$ ; M is selected from  $-C(=O)N(R^7)-$  and  $-SO_2N(R^7)-$ ; and any C-H of  $L^2$  is optionally  $C-R^{20}$ ;

ring C is either a five- to ten-membered aryl or a five- to ten-membered heteroaryl; each  $R^3$  is independently selected from -H, halogen, trihalomethyl, -CN,  $-NO_2$ ,  $-OR^4$ ,  $-N(R^4)R^4$ ,  $-S(O)_{0-2}R^4$ ,  $-SO_2N(R^4)R^4$ ,  $-CO_2R^4$ ,  $-C(=O)N(R^4)R^4$ ,  $-C(=NR^5)N(R^4)R^4$ ,  $-C(=NR^5)R^4$ ,  $-N(R^4)SO_2R^4$ ,  $-N(R^4)C(O)R^4$ ,  $-NCO_2R^4$ ,  $-C(=O)R^4$ , optionally substituted alkoxy, optionally substituted  $C_{1-6}$ alkyl, optionally substituted aryl, optionally substituted aryl  $C_{1-6}$ alkyl, optionally substituted heterocyclyl, and optionally substituted heterocyclyl  $C_{1-6}$ alkyl; provided  $R^3$  is not a cyclic sulfonamide attached to ring C via the nitrogen of said cyclic sulfonamide;

two adjacent of  $R^3$ , together with the annular atoms to which they are attached, can form a five- to six-membered ring containing up to two heteroatoms and optionally substituted with up to three of  $R^{25}$ ;

$R^4$  is selected from -H, optionally substituted  $C_{1-6}$ alkyl, optionally substituted aryl, optionally substituted aryl  $C_{1-6}$ alkyl, optionally substituted heterocyclyl, and optionally substituted heterocyclyl  $C_{1-6}$ alkyl;

two of  $R^4$ , when taken together with a common nitrogen to which they are attached, form an optionally substituted five- to seven-membered heterocyclyl, said optionally substituted five- to seven-membered heterocyclyl optionally containing at least one additional heteroatom selected from N, O, S, and P;

$R^5$  is selected from -H, -CN,  $-NO_2$ ,  $-OR^4$ ,  $-S(O)_{0-2}R^4$ ,  $-CO_2R^4$ , optionally substituted  $C_{1-6}$ alkyl, optionally substituted  $C_{1-6}$ alkenyl, and optionally substituted  $C_{1-6}$ alkynyl;

$R^7$  is selected from -H, optionally substituted  $C_{1-6}$ alkyl,  $-SO_2N(R^4)R^4$ ,  $-CO_2R^4$ ,  $-C(=O)N(R^4)R^4$ ,  $-C(=NR^5)N(R^4)R^4$ ,  $-C(=NR^5)R^4$ ,  $-C(=O)R^4$ , optionally substituted alkoxy, optionally substituted aryl, optionally substituted aryl  $C_{1-6}$ alkyl, optionally substituted heterocyclyl, and optionally substituted heterocyclyl  $C_{1-6}$ alkyl; and

each of R<sup>10</sup>, each of R<sup>15</sup>, each of R<sup>20</sup>, and each of R<sup>25</sup> is independently selected from -H, halogen, trihalomethyl, -CN, -NO<sub>2</sub>, -OR<sup>4</sup>, -N(R<sup>4</sup>)R<sup>4</sup>, -S(O)<sub>0-2</sub>R<sup>4</sup>, -SO<sub>2</sub>N(R<sup>4</sup>)R<sup>4</sup>, -CO<sub>2</sub>R<sup>4</sup>, -C(=O)N(R<sup>4</sup>)R<sup>4</sup>, -C(=NR<sup>5</sup>)N(R<sup>4</sup>)R<sup>4</sup>, -C(=NR<sup>5</sup>)R<sup>4</sup>, -N(R<sup>4</sup>)SO<sub>2</sub>R<sup>4</sup>, -N(R<sup>4</sup>)C(O)R<sup>4</sup>, -NCO<sub>2</sub>R<sup>4</sup>, -C(=O)R<sup>4</sup>, optionally substituted alkoxy, optionally substituted C<sub>1-6</sub>alkyl, optionally substituted aryl, optionally substituted aryl C<sub>1-6</sub>alkyl, optionally substituted heterocyclyl, and optionally substituted heterocyclyl C<sub>1-6</sub>alkyl;

provided:

1) when both ring B and ring C are phenyl:

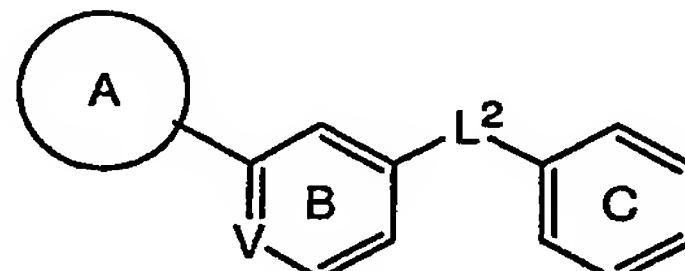
- a) and the compound comprises ring B-CH<sub>2</sub>N(H)C(=O)N(H)-ring C, then L<sup>1</sup> must be a single bond; R<sup>3</sup> can not comprise a group of the formula -O(CH<sub>2</sub>)<sub>2-4</sub>-N-piperazine that is *ortho*- to L<sup>2</sup>; and ring A cannot be a 5-methyl-[1,2,4]-oxadiazol-3-yl radical, a 4H-[1,2,4]-oxadiazol-5-one-3-yl radical, nor a 4'-[2,2';6',2"]terpyridinyl radical;
- b) and L<sup>1</sup> is single bond, then L<sup>2</sup> cannot comprise -N(H)C(=O)C(=O)N(H)- nor -N(H)C(=Q)C(H)CNC(=O)- (where Q is S or O);
- c) and L<sup>1</sup> is other than single bond, then A cannot be quinolin-2-yl-L<sup>1</sup>, quinolin-3-yl-L<sup>1</sup>, or quinolin-4-yl-L<sup>1</sup>;

2) when ring A is a fused aryl system, then L<sup>1</sup> must be a single bond;

3) when ring B is phenyl, ring C is a C<sub>6-16</sub>carbocyclic, L<sup>1</sup> is a single bond, and the compound comprises -ring B-OCH<sub>2</sub>C(=O)N(H)- then ring A cannot be a 2,5-dimethyl-1H-pyrrole-1-yl radical;

4) ring A cannot be a pyrimidin-2-yl radical when L<sup>1</sup> is -N(H)- and ring B is phenyl;

5) when the compound comprises the formula,

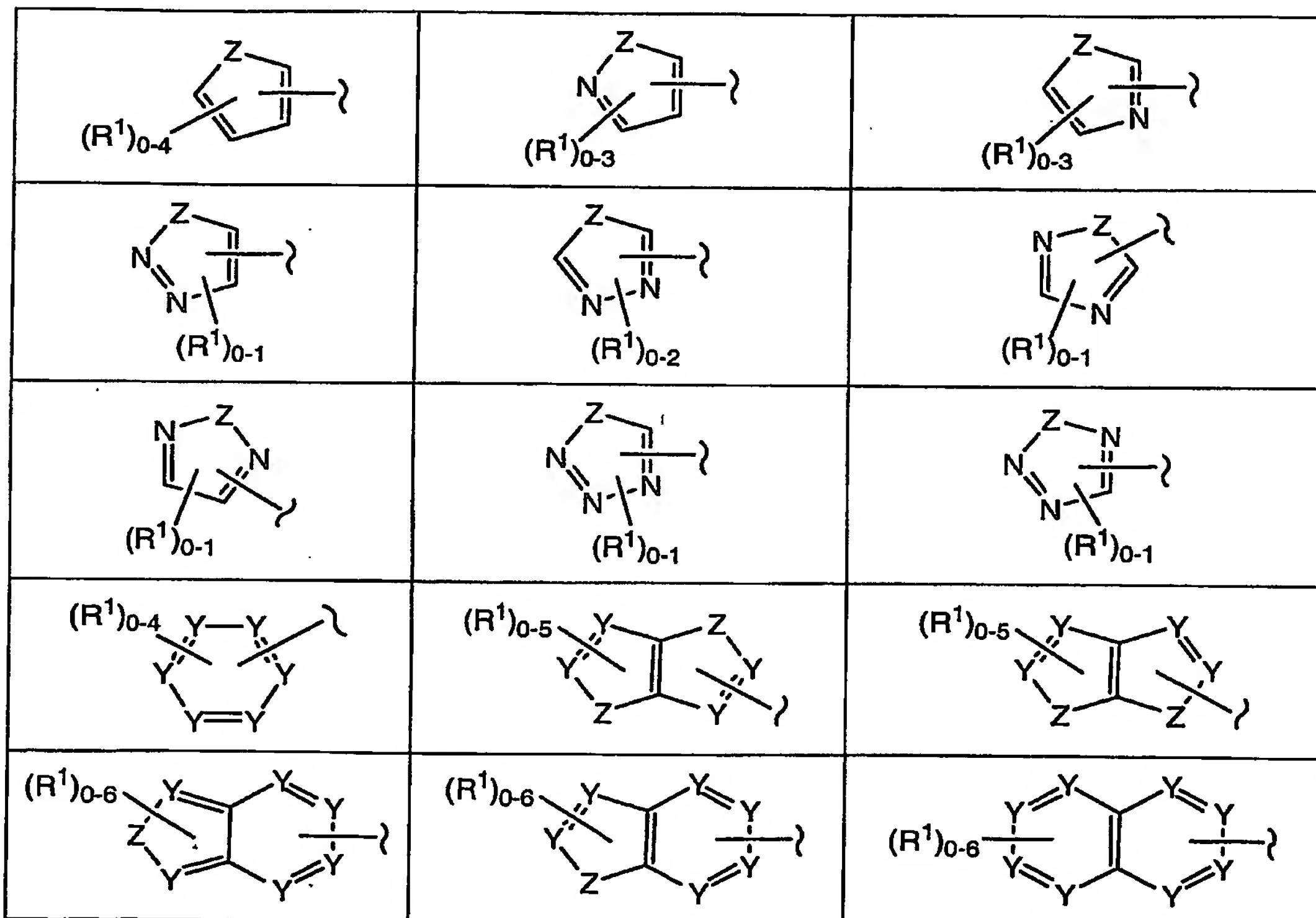


where V is =C(H)- or =N-, and there is a nitrogen of L<sup>2</sup> bound directly to ring B, then A can not comprise a [1,2,4]-oxadiazol-3-yl radical; and

- 6) the compound is not one of: N-naphthalen-1-yl-2-[{3-(1H-tetrazol-1-yl)phenyl]oxy}acetamide, N-[4-(phenyloxy)phenyl]-2-[{3-(1H-tetrazol-1-yl)phenyl]oxy}acetamide, N-(3,4-dimethylphenyl)-2-[{3-(1H-tetrazol-1-yl)phenyl]oxy}acetamide, N-(2,3-dimethylphenyl)-2-[{3-(1H-tetrazol-1-yl)phenyl]oxy}acetamide, N-(2,4-dimethylphenyl)-2-[{3-(1H-tetrazol-1-yl)phenyl]oxy}acetamide, N-(2,5-dimethylphenyl)-2-[{3-(1H-tetrazol-1-yl)phenyl]oxy}acetamide, N-(3,5-dimethylphenyl)-2-[{3-(1H-tetrazol-1-yl)phenyl]oxy}acetamide, N-(2,6-dimethylphenyl)-2-[{3-(1H-tetrazol-1-yl)phenyl]oxy}acetamide, 2-[{3-(1H-tetrazol-1-yl)phenyl]oxy}-N-(2,4,6-trimethylphenyl)acetamide, N-(2-ethylphenyl)-2-[{3-(1H-tetrazol-1-yl)phenyl]oxy}acetamide, N-(4-ethylphenyl)-2-[{3-(1H-tetrazol-1-yl)phenyl]oxy}acetamide, N-(2,6-diethylphenyl)-2-[{3-(1H-tetrazol-1-yl)phenyl]oxy}acetamide, N-[2-(methyloxy)phenyl]-2-[{3-(1H-tetrazol-1-yl)phenyl]oxy}acetamide, N-[2-(ethyloxy)phenyl]-2-[{3-(1H-tetrazol-1-yl)phenyl]oxy}acetamide, N-[3-(ethyloxy)phenyl]-2-[{3-(1H-tetrazol-1-yl)phenyl]oxy}acetamide, N-[2,4-bis(methyloxy)phenyl]-2-[{3-(1H-tetrazol-1-yl)phenyl]oxy}acetamide, N-[4-(dimethylamino)phenyl]-2-[{3-(1H-tetrazol-1-yl)phenyl]oxy}acetamide, N-(2,3-dichlorophenyl)-2-[{3-(1H-tetrazol-1-yl)phenyl]oxy}acetamide, N-(4-chloro-3-methylphenyl)-2-[{3-(1H-tetrazol-1-yl)phenyl]oxy}acetamide, N-(4-bromophenyl)-2-[{3-(1H-tetrazol-1-yl)phenyl]oxy}acetamide, N-(2-fluorophenyl)-2-[{3-(1H-tetrazol-1-yl)phenyl]oxy}acetamide, N-(4-fluorophenyl)-2-[{3-(1H-tetrazol-1-yl)phenyl]oxy}acetamide, 2-[{3-(1H-tetrazol-1-yl)phenyl]oxy}-N-[2-(trifluoromethyl)phenyl]acetamide, 2-[{3-(1H-tetrazol-1-yl)phenyl]oxy}-N-[3-(trifluoromethyl)phenyl]acetamide, methyl 4-[({[3-(1H-tetrazol-1-yl)phenyl]oxy}acetyl)amino]benzoate, ethyl 4-[({[3-(1H-tetrazol-1-yl)phenyl]oxy}acetyl)amino]benzoate, 3-[({[3-(1H-tetrazol-1-yl)phenyl]oxy}acetyl)amino]benzoic acid, N-[3-(methyloxy)phenyl]-2-[{3-(1H-tetrazol-1-yl)phenyl]oxy}acetamide, N-[4-(methyloxy)phenyl]-2-[{3-(1H-tetrazol-1-yl)phenyl]oxy}acetamide, N-[2-chloro-5-(trifluoromethyl)phenyl]-2-[{3-(1H-tetrazol-1-yl)phenyl]oxy}acetamide, N-[4-chloro-3-(trifluoromethyl)phenyl]-2-[{3-(4H-1,2,4-triazol-4-yl)phenyl]oxy}acet-

amide, N-(4-chlorophenyl)-2-{[3-(1H-tetrazol-1-yl)phenyl]oxy}acetamide, N-(4-aminophenyl)-2-{[3-(1H-tetrazol-1-yl)phenyl]oxy}acetamide, and N-(4-acetylphenyl)-2-{[3-(1H-tetrazol-1-yl)phenyl]oxy}acetamide.

2. The compound according to claim 1, wherein L<sup>1</sup> is a single bond.
3. The compound according to claim 2, wherein ring A contains between one and four annular nitrogens.
4. The compound according to claim 3, wherein ring A is selected from the following:



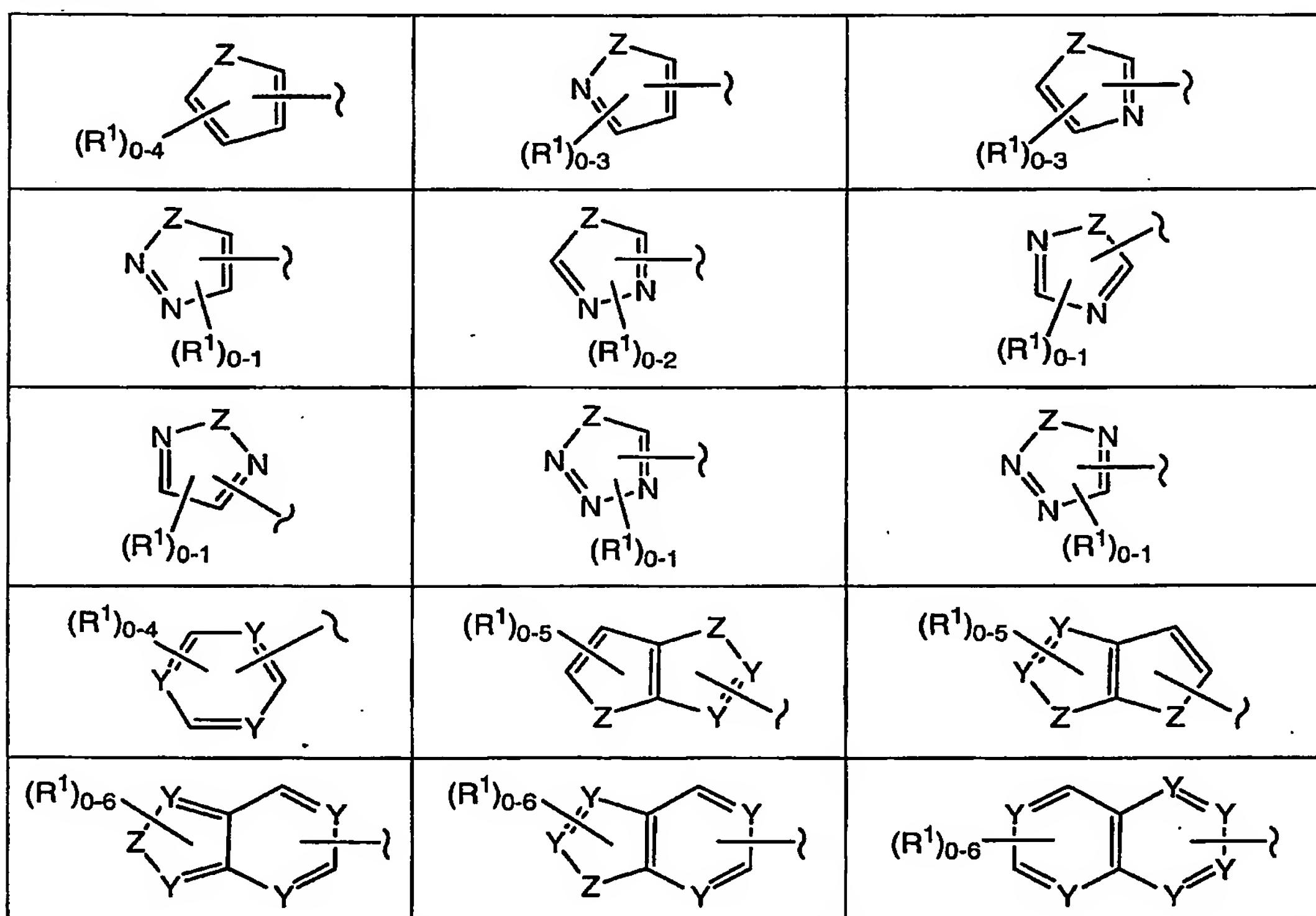
wherein each Y is independently either =C(H)- or =N-; and Z is selected from -O-, -S(O)<sub>0-2</sub>-, and -N(R<sup>7</sup>)-.

5. The compound according to claim 4, wherein ring B is phenylene or pyridylene.
6. The compound according to claim 5, wherein the annular atoms of ring B to which L<sup>1</sup> and L<sup>2</sup> are attached are not contiguous.

7. The compound according to claim 6, wherein  $L^2$  is selected from  $-X(CH_2)_2O-$ ,  $-X(CH_2)_2N(R^7)-$ ,  $-CH_2XC(=O)N(R^7)-$ ,  $-XCH_2SO_2N(R^7)-$ ,  $-XN(R^7)C(=O)N(R^7)-$  and  $-XCH_2C(=O)N(R^7)-$ ; wherein X is selected from  $-CH_2-$ ,  $-O-$ ,  $-S(O)_{0-2}-$  and  $-N(R^7)-$ ; and any C-H of  $L^2$  is optionally  $C-R^{20}$ .

8. The compound according to claim 7, wherein  $L^2$  is selected from  $-N(H)N(H)C(=O)N(H)-$ ,  $-CH_2N(H)C(=O)N(H)-$ ,  $-CH_2OC(=O)N(H)-$ , and  $-XCH_2C(=O)N(H)-$ ; wherein X is selected from  $-O-$ ,  $-S(O)_{0-2}-$ , and  $-N(R^7)-$ ; and any C-H of  $L^2$  is optionally  $C-R^{20}$ .

9. The compound according to claim 8, wherein ring A is selected from the following:



wherein each Y is independently either  $=C(H)-$  or  $=N-$ ; and Z is selected from  $-O-$ ,  $-S-$ , and  $-N(R^7)-$ .

10. The compound according to claim 9, wherein ring C is phenyl or pyridyl.

11. The compound according to claim 10, wherein there exists at least one of  $R^3$  that is halogen.

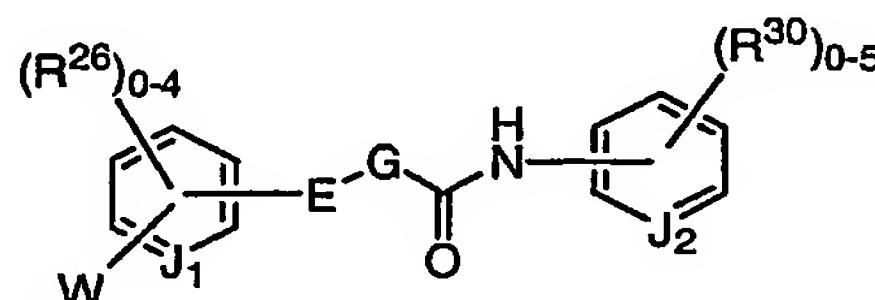
12. The compound according to claim 10, wherein there exists at least one of R<sup>3</sup> that is trihalomethyl.

13. The compound according to claim 10, wherein there exists at least one of R<sup>3</sup> that is trifluoromethyl.

14. The compound according to claim 13, wherein ring C is a phenyl comprising a trifluoromethyl radical *meta*- to L<sup>2</sup>.

15. The compound according to claim 10, wherein each of R<sup>3</sup> is independently selected from -H, halogen, trihalomethyl, -OR<sup>4</sup>, -CO<sub>2</sub>R<sup>4</sup>, -C(=O)R<sup>4</sup>, and optionally substituted C<sub>1</sub>-alkyl.

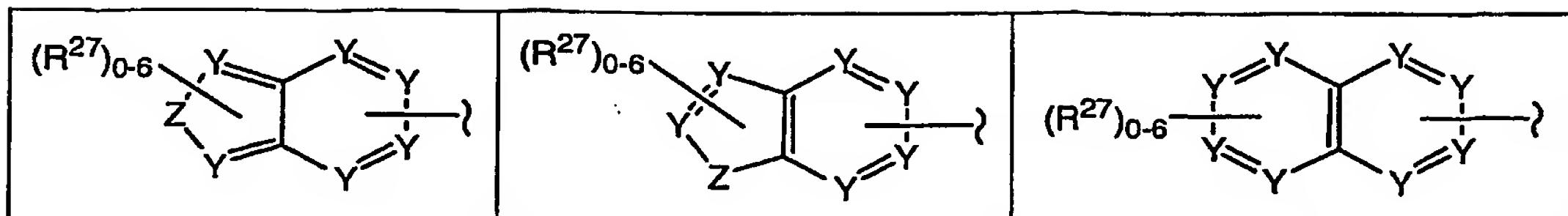
16. A compound for modulating c-Kit activity according to Formula III,



III

or a pharmaceutically acceptable salt, hydrate, or prodrug thereof, wherein,

W is selected from the following:

each of R<sup>27</sup> independently selected from halogen, trihalomethyl, -CN, -NO<sub>2</sub>, -OR<sup>55</sup>, -N(R<sup>55</sup>)R<sup>55</sup>, -S(O)<sub>0-2</sub>R<sup>55</sup>, -SO<sub>2</sub>N(R<sup>55</sup>)R<sup>55</sup>, -CO<sub>2</sub>R<sup>55</sup>, -C(=O)N(R<sup>55</sup>)R<sup>55</sup>, -C(=NR<sup>50</sup>)N(R<sup>55</sup>)R<sup>55</sup>, -C(=NR<sup>50</sup>)R<sup>55</sup>, -N(R<sup>55</sup>)SO<sub>2</sub>R<sup>55</sup>, -N(R<sup>55</sup>)C(O)R<sup>55</sup>, -NCO<sub>2</sub>R<sup>55</sup>, -C(=O)R<sup>55</sup>, optionally substituted alkoxy, optionally substituted C<sub>1-6</sub>alkyl, optionally substituted aryl, optionally substituted aryl C<sub>1-6</sub>alkyl, optionally substituted heterocyclyl, and optionally substituted heterocyclyl C<sub>1-6</sub>alkyl;

each Y is independently either =C(H)- or =N-;

Z is selected from -O-, -S(O)<sub>0-2</sub>-, and -N(R<sup>7</sup>)-

E and G are each independently selected from -O-, -S(O)<sub>0-2</sub>-, -C(R<sup>31</sup>)R<sup>32</sup>-, and -N(R<sup>33</sup>)-;

J<sub>1</sub> and J<sub>2</sub> are each independently =C(H)- or =N-;

each of R<sup>26</sup> and R<sup>30</sup> is independently selected from -H, halogen, trihalomethyl, -CN, -NO<sub>2</sub>, -OR<sup>40</sup>, -N(R<sup>40</sup>)R<sup>40</sup>, -S(O)<sub>0-2</sub>R<sup>40</sup>, -SO<sub>2</sub>N(R<sup>40</sup>)R<sup>40</sup>, -CO<sub>2</sub>R<sup>40</sup>, -C(=O)N(R<sup>40</sup>)R<sup>40</sup>, -C(=NR<sup>50</sup>)N(R<sup>40</sup>)R<sup>40</sup>, -C(=NR<sup>50</sup>)R<sup>40</sup>, -N(R<sup>40</sup>)SO<sub>2</sub>R<sup>40</sup>, -N(R<sup>40</sup>)C(O)R<sup>40</sup>, -NCO<sub>2</sub>R<sup>40</sup>, -C(=O)R<sup>40</sup>, optionally substituted alkoxy, optionally substituted C<sub>1-6</sub>alkyl, optionally substituted aryl, optionally substituted aryl C<sub>1-6</sub>alkyl, optionally substituted heterocyclyl, and optionally substituted heterocyclyl C<sub>1-6</sub>alkyl;

two adjacent of R<sup>26</sup> or two adjacent of R<sup>30</sup>, together with the annular atoms to which they are attached, can form a five- to six-membered ring containing up to two heteroatoms and optionally substituted with up to three of R<sup>35</sup>;

R<sup>31</sup> and R<sup>32</sup> are each independently selected from -H, halogen, trihalomethyl, -CN, -NO<sub>2</sub>, -OR<sup>40</sup>, -N(R<sup>40</sup>)R<sup>40</sup>, -S(O)<sub>0-2</sub>R<sup>40</sup>, -SO<sub>2</sub>N(R<sup>40</sup>)R<sup>40</sup>, -CO<sub>2</sub>R<sup>40</sup>, -C(=O)N(R<sup>40</sup>)R<sup>40</sup>, -C(=NR<sup>50</sup>)N(R<sup>40</sup>)R<sup>40</sup>, -C(=NR<sup>50</sup>)R<sup>40</sup>, -N(R<sup>40</sup>)SO<sub>2</sub>R<sup>40</sup>, -N(R<sup>40</sup>)C(O)R<sup>40</sup>, -NCO<sub>2</sub>R<sup>40</sup>, -C(=O)R<sup>40</sup>, optionally substituted alkoxy, optionally substituted C<sub>1-6</sub>alkyl, optionally substituted aryl, optionally substituted aryl C<sub>1-6</sub>alkyl, optionally substituted heterocyclyl, and optionally substituted heterocyclyl C<sub>1-6</sub>alkyl;

R<sup>33</sup> is selected from -H, optionally substituted lower alkyl, -SO<sub>2</sub>N(R<sup>40</sup>)R<sup>40</sup>, -CO<sub>2</sub>R<sup>40</sup>, -C(=O)N(R<sup>40</sup>)R<sup>40</sup>, -C(=NR<sup>50</sup>)N(R<sup>40</sup>)R<sup>40</sup>, -C(=NR<sup>50</sup>)R<sup>40</sup>, -C(=O)R<sup>40</sup>, optionally

substituted alkoxy, optionally substituted C<sub>1-6</sub>alkyl, optionally substituted aryl, optionally substituted aryl C<sub>1-6</sub>alkyl, optionally substituted heterocyclyl, and optionally substituted heterocyclyl C<sub>1-6</sub>alkyl;

R<sup>40</sup> is selected from -H, optionally substituted alkoxy, optionally substituted C<sub>1-6</sub>alkyl, optionally substituted aryl, optionally substituted aryl C<sub>1-6</sub>alkyl, optionally substituted heterocyclyl, and optionally substituted heterocyclyl C<sub>1-6</sub>alkyl;

two of R<sup>40</sup>, when taken together with a common nitrogen to which they are attached, form an optionally substituted five- to seven-membered heterocyclyl, said optionally substituted five- to seven-membered heterocyclyl optionally containing at least one additional heteroatom selected from N, O, S, and P;

R<sup>50</sup> is selected from -H, -CN, -NO<sub>2</sub>, -OR<sup>40</sup>, -S(O)<sub>0-2</sub>R<sup>40</sup>, -CO<sub>2</sub>R<sup>40</sup>, optionally substituted C<sub>1-6</sub>alkyl, optionally substituted C<sub>1-6</sub>alkenyl, and optionally substituted C<sub>1-6</sub>alkynyl;

R<sup>55</sup> is selected from -H, optionally substituted C<sub>1-6</sub>alkyl, optionally substituted heterocyclyl, and optionally substituted heterocyclyl C<sub>1-6</sub>alkyl; and

two of R<sup>55</sup>, when taken together with a common nitrogen to which they are attached, form an optionally substituted five- to seven-membered heterocyclyl, said optionally substituted five- to seven-membered heterocyclyl optionally containing at least one additional heteroatom selected from N, O, S, and P.

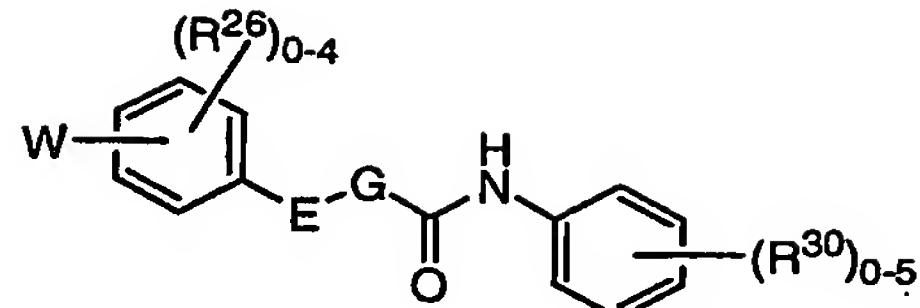
17. The compound according to claim 16, wherein the annular carbons of ring B to which W and E are attached are not contiguous.

18. The compound according to claim 17, wherein R<sup>30</sup> is selected from -H, halogen, trihalomethyl, -OR<sup>40</sup>, -N(R<sup>40</sup>)R<sup>40</sup>, -CO<sub>2</sub>R<sup>40</sup>, -C(=O)R<sup>40</sup>, optionally substituted alkoxy, optionally substituted C<sub>1-6</sub>alkyl, optionally substituted aryl, optionally substituted aryl C<sub>1-6</sub>alkyl, optionally substituted heterocyclyl, and optionally substituted heterocyclyl C<sub>1-6</sub>alkyl.

19. The compound according to claim 18, wherein there exists at least one of R<sup>30</sup> that is trihalomethyl.

20. The compound according to claim 18, wherein there exists at least one of R<sup>30</sup> that is trifluoromethyl.

21. The compound according to claim 18, according to formula III.



22. The compound according to claim 21, wherein W is selected from the following:


and R<sup>27</sup> is defined as above.

23. The compound according to claim 22, wherein E is selected from -O-, -S(O)<sub>0-2</sub>-, and -NH-; and G is -CH<sub>2</sub>-.

24. The compound according to claim 22, wherein E is either -CH<sub>2</sub>- or -NH-; and G is selected from -O-, -S-, and -NH-.

25. The compound according to either claim 23 or claim 24, wherein each of R<sup>3</sup> is independently selected from -H, halogen, trihalomethyl, -OR<sup>4</sup>, -CO<sub>2</sub>R<sup>4</sup>, -C(=O)R<sup>4</sup>, and optionally substituted C<sub>1-6</sub>alkyl.

26. The compound according to claim 25, wherein at least one of R<sup>30</sup> is a trifluoromethyl radical *meta*- to -E-G-C(=O)N(H)-.

27. The compound according to either claim 1 or claim 16, selected from Table 3:

Table 3

Entry	Name	Structure
1	N-[5-chloro-2-(methyloxy)phenyl]-2-{[3-(1H-tetrazol-1-yl)phenyl]oxy}acetamide	
2	N-phenyl-2-{[3-(1H-tetrazol-1-yl)phenyl]oxy}acetamide	
3	N-(2-methylphenyl)-2-{[3-(1H-tetrazol-1-yl)phenyl]oxy}acetamide	
4	N-(2-chlorophenyl)-2-{[3-(1H-tetrazol-1-yl)phenyl]oxy}acetamide	
5	N-[4-chloro-3-(trifluoromethyl)phenyl]-2-{[3-(1H-tetrazol-1-yl)phenyl]oxy}acetamide	
6	ethyl 2-[( {[3-(1H-tetrazol-1-yl)phenyl]oxy}acetyl)amino]-4,5,6,7-tetrahydro-1-benzothiophene-3-carboxylate	

Table 3

Entry	Name	Structure
7	N-(3-chloro-2-methylphenyl)-2-{[3-(1H-tetrazol-1-yl)phenyl]oxy}acetamide	
8	N-(3-fluorophenyl)-2-{[3-(1H-tetrazol-1-yl)phenyl]oxy}acetamide	
9	N-[4-chloro-3-(trifluoromethyl)phenyl]-2-{[3-(2H-tetrazol-5-yl)phenyl]oxy}acetamide	
10	N-(4-chloro-2-fluorophenyl)-2-{[3-(1H-tetrazol-1-yl)phenyl]oxy}acetamide	
11	N-(4-bromo-3-methylphenyl)-2-{[3-(1H-tetrazol-1-yl)phenyl]oxy}acetamide	
12	N-(4-morpholin-4-ylphenyl)-2-{[3-(1H-tetrazol-1-yl)phenyl]oxy}acetamide	
13	N-[4-fluoro-3-(trifluoromethyl)phenyl]-2-{[3-(1H-tetrazol-1-yl)phenyl]oxy}acetamide	

Table 3

Entry	Name	Structure
14	N-[4-bromo-3-(trifluoromethyl)phenyl]-2-{[3-(1H-tetrazol-1-yl)phenyl]oxy}acetamide	
15	N-[4-chloro-3-(trifluoromethyl)phenyl]-2-{[4-(1H-tetrazol-1-yl)phenyl]oxy}acetamide	
16	N-[4-chloro-3-(trifluoromethyl)phenyl]-2-{[3-(1H-tetrazol-1-yl)phenyl]oxy}propanamide	
17	N-[4-chloro-3-(trifluoromethyl)phenyl]-2-{[3-(5-methyl-1H-tetrazol-1-yl)phenyl]oxy}acetamide	
18	N-[4-chloro-3-(trifluoromethyl)phenyl]-2-{[2-methyl-5-(1H-tetrazol-1-yl)phenyl]oxy}acetamide	
19	N-(4-chlorophenyl)-N-methyl-2-{[3-(1H-tetrazol-1-yl)phenyl]oxy}acetamide	
20	N-[4-chloro-2-(trifluoromethyl)phenyl]-2-{[3-(1H-tetrazol-1-yl)phenyl]oxy}acetamide	

Table 3

Entry	Name	Structure
21	N-[4-chloro-3-(trifluoromethyl)phenyl]-2-{[3-(2,5-dioxopyrrolidin-1-yl)phenyl]oxy}acetamide	
22	(2E)-N-[4-chloro-3-(trifluoromethyl)phenyl]-3-[3-(1H-tetrazol-1-yl)phenyl]prop-2-enamide	
23	N-[4-fluoro-3-(trifluoromethyl)phenyl]-2-{[4-(1H-tetrazol-1-yl)phenyl]oxy}acetamide	
24	N-[4-chloro-3-(trifluoromethyl)phenyl]-2-{[3-(2-methyl-2H-tetrazol-5-yl)phenyl]oxy}acetamide	
25	N-[4-chloro-3-(trifluoromethyl)phenyl]-2-{[2,4-dichloro-5-(1H-tetrazol-1-yl)phenyl]oxy}acetamide	
26	N-[4-chloro-3-(trifluoromethyl)phenyl]-2-{[3-(1H-tetrazol-1-yl)phenyl]thio}acetamide	
27	N-[4-chloro-3-(trifluoromethyl)phenyl]-N~2-[3-(1H-tetrazol-1-yl)phenyl]glycinamide	

Table 3

Entry	Name	Structure
28	N-[4-chloro-3-(trifluoromethyl)phenyl]-2-{[2-(1H-tetrazol-1-yl)phenyl]oxy}acetamide	
29	methyl 1-{3-[(2-{[4-chloro-3-(trifluoromethyl)phenyl]amino}-2-oxoethyl)oxy]phenyl}-1H-1,2,3-triazole-4-carboxylate	
30	1,1-dimethylethyl {4-[{[3-(1H-tetrazol-1-yl)phenyl]oxy}acetyl]amino}phenyl carbamate	
31	1,1-dimethylethyl {4-[{[4-(1H-tetrazol-1-yl)phenyl]oxy}acetyl]amino}phenyl carbamate	
32	N-{4-[(1-ethylpiperidin-4-yl)amino]phenyl}-2-{[3-(1H-tetrazol-1-yl)phenyl]oxy}acetamide	
33	N-{4-[(1-ethylpiperidin-3-yl)amino]phenyl}-2-{[3-(1H-tetrazol-1-yl)phenyl]oxy}acetamide	
34	N-(4-aminophenyl)-2-{[4-(1H-tetrazol-1-yl)phenyl]oxy}acetamide	

Table 3

Entry	Name	Structure
35	N-{4-[(1-ethylpiperidin-4-yl)amino]phenyl}-2-{[4-(1H-tetrazol-1-yl)phenyl]oxy}acetamide	
36	N-{4-[(1-ethylpiperidin-3-yl)amino]phenyl}-2-{[4-(1H-tetrazol-1-yl)phenyl]oxy}acetamide	
37	N-[4-chloro-3-(trifluoromethyl)phenyl]-2-[(3-pyridin-4-ylphenyl)oxy]acetamide	
38	N-[4-chloro-3-(trifluoromethyl)phenyl]-N~2~-methyl-N~2~-[3-(1H-tetrazol-1-yl)phenyl]glycinamide	
39	N-1,3-benzothiazol-2-yl-2-{[3-(1H-tetrazol-1-yl)phenyl]oxy}acetamide	
40	N-quinolin-8-yl-2-{[3-(1H-tetrazol-1-yl)phenyl]oxy}acetamide	
41	N-(2,3-dihydro-1,4-benzodioxin-6-yl)-2-{[3-(1H-tetrazol-1-yl)phenyl]oxy}acetamide	

Table 3

Entry	Name	Structure
42	N-isoquinolin-5-yl-2-{[3-(1H-tetrazol-1-yl)phenyl]oxy}acetamide	
43	N-{3-[(phenylmethyl)oxy]phenyl}-2-{[3-(1H-tetrazol-1-yl)phenyl]oxy}acetamide	
44	N-[5-methyl-2-(methyloxy)phenyl]-2-{[3-(1H-tetrazol-1-yl)phenyl]oxy}acetamide	
45	N-[2,5-bis(methyloxy)phenyl]-2-{[3-(1H-tetrazol-1-yl)phenyl]oxy}acetamide	
46	N-(6-fluoro-1,3-benzothiazol-2-yl)-2-{[3-(1H-tetrazol-1-yl)phenyl]oxy}acetamide	
47	methyl 3-[(2-{[3-(1H-tetrazol-1-yl)phenyl]oxy}acetyl)amino]benzoate	
48	5-chloro-2-[(2-{[3-(1H-tetrazol-1-yl)phenyl]oxy}acetyl)amino]benzamide	

Table 3

Entry	Name	Structure
49	N-[5-chloro-2,4-bis(methyloxy)phenyl]-2-{[3-(1H-tetrazol-1-yl)phenyl]oxy}acetamide	
50	N-[2-(phenyloxy)phenyl]-2-{[3-(1H-tetrazol-1-yl)phenyl]oxy}acetamide	
51	N-[3-(aminosulfonyl)phenyl]-2-{[3-(1H-tetrazol-1-yl)phenyl]oxy}acetamide	
52	N-[2-(methyloxy)-5-(trifluoromethyl)phenyl]-2-{[3-(1H-tetrazol-1-yl)phenyl]oxy}acetamide	
53	N-(4-[(4-methylphenyl)sulfonyl]amino)phenyl)-2-{[3-(1H-tetrazol-1-yl)phenyl]oxy}acetamide	
54	N-(5-phenyl-1H-pyrazol-3-yl)-2-{[3-(1H-tetrazol-1-yl)phenyl]oxy}acetamide	
55	N-1,3-benzothiazol-2-yl-2-{[4-(1H-tetrazol-1-yl)phenyl]oxy}acetamide	

Table 3

Entry	Name	Structure
56	N-quinolin-8-yl-2-{{4-(1H-tetrazol-1-yl)phenyl}oxy}acetamide	
57	1,1-dimethylethyl 2-{3-[{2-{{4-chloro-3-(trifluoromethyl)phenyl}amino}-2-oxoethyl}oxy]phenyl}-1H-pyrrole-1-carboxylate	
58	N-[4-chloro-3-(trifluoromethyl)phenyl]-2-{{3-(1H-pyrrol-2-yl)phenyl}oxy}acetamide	
59	N-[4-chloro-3-(trifluoromethyl)phenyl]-2-{{3-(pyrimidin-5-yl)phenyl}oxy}acetamide	
60	N-[4-chloro-3-(trifluoromethyl)phenyl]-2-{{3-(1H-1,2,3-triazol-1-yl)phenyl}oxy}acetamide	
61	4-chloro-N-(2-{{3-(1H-tetrazol-1-yl)phenyl}oxy}ethyl)-3-(trifluoromethyl)aniline	
62	N-[4-chloro-3-(trifluoromethyl)phenyl]-N-(2-{{3-(1H-tetrazol-1-yl)phenyl}oxy}ethyl)formamide	

Table 3

Entry	Name	Structure
63	N-[4-chloro-3-(trifluoromethyl)phenyl]-2-[(3-pyridin-3-ylphenyl)oxy]acetamide	
64	N-[4-chloro-3-(trifluoromethyl)phenyl]-2-[(3-furan-3-ylphenyl)oxy]acetamide	
65	(2E)-N-[4-fluoro-3-(trifluoromethyl)phenyl]-3-[3-(1H-tetrazol-1-yl)phenyl]prop-2-enamide	
66	N-[4-fluoro-3-(trifluoromethyl)phenyl]-3-[3-(1H-tetrazol-1-yl)phenyl]propanamide	
67	N-[4-chloro-3-(trifluoromethyl)phenyl]-2-[(6-(1H-tetrazol-1-yl)pyrimidin-4-yl)oxy]acetamide	
68	N-[4-chloro-3-(trifluoromethyl)phenyl]-2-[(3-(3,5-dimethylisoxazol-4-yl)phenyl)oxy]acetamide	
69	N-[4-chloro-3-(trifluoromethyl)phenyl]-2-[(3-quinolin-7-ylphenyl)oxy]acetamide	

Table 3

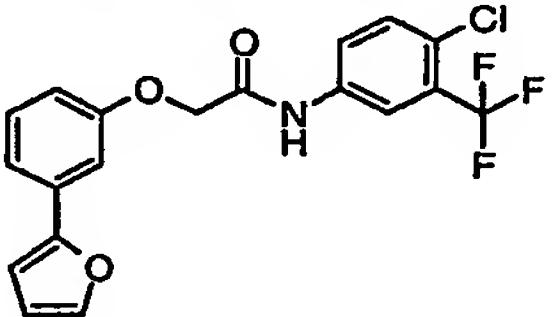
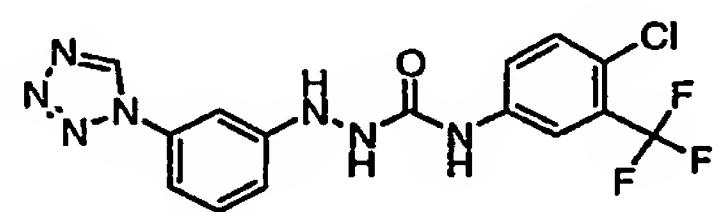
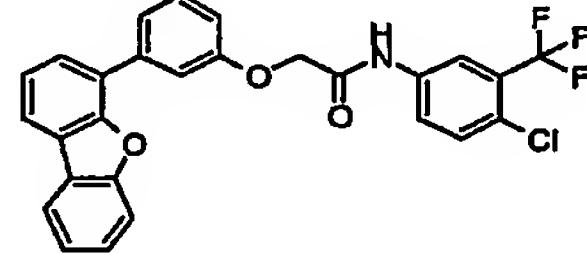
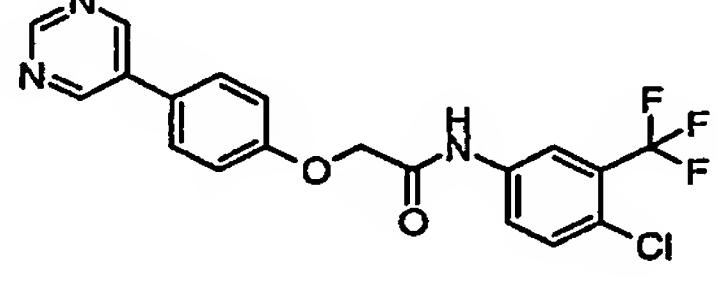
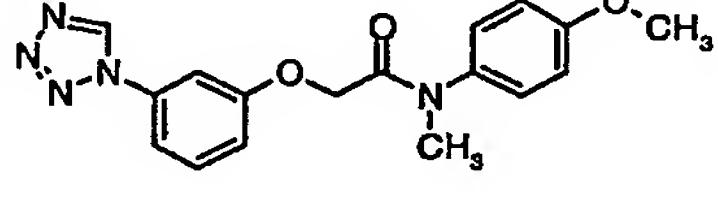
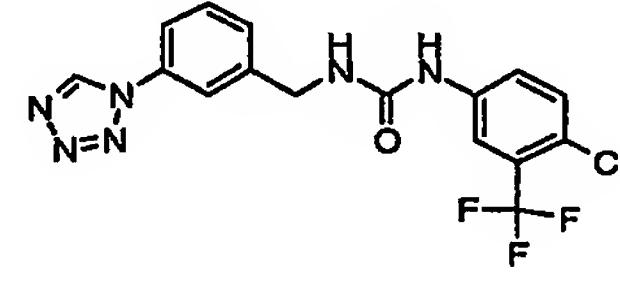
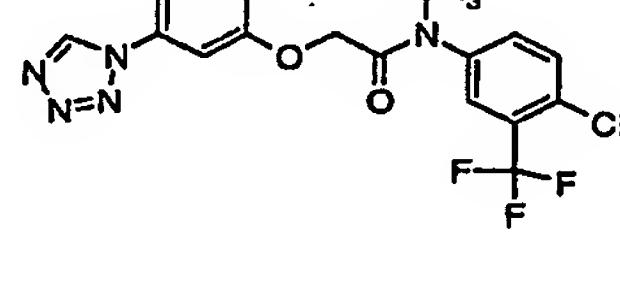
Entry	Name	Structure
70	N-[4-chloro-3-(trifluoromethyl)phenyl]-2-[(3-furan-2-ylphenyl)oxy]acetamide	
71	N-[4-chloro-3-(trifluoromethyl)phenyl]-2-[3-(1H-tetrazol-1-yl)phenyl]hydrazinecarboxamide	
72	N-[4-chloro-3-(trifluoromethyl)phenyl]-2-[(3-dibenzo[b,d]furan-4-ylphenyl)oxy]acetamide	
73	N-[4-chloro-3-(trifluoromethyl)phenyl]-2-[(4-pyrimidin-5-ylphenyl)oxy]acetamide	
74	N-methyl-N-[4-(methoxy)phenyl]-2-{[3-(1H-tetrazol-1-yl)phenyl]oxy}acetamide	
75	N-[4-chloro-3-(trifluoromethyl)phenyl]-N'-{[3-(1H-tetrazol-1-yl)phenyl]methyl}urea	
76	N-[4-chloro-3-(trifluoromethyl)phenyl]-N-methyl-2-{[3-(1H-tetrazol-1-yl)phenyl]oxy}acetamide	

Table 3

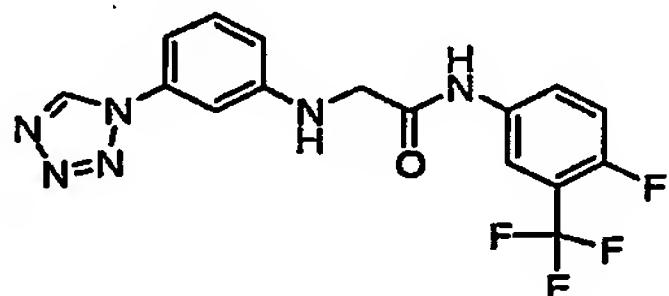
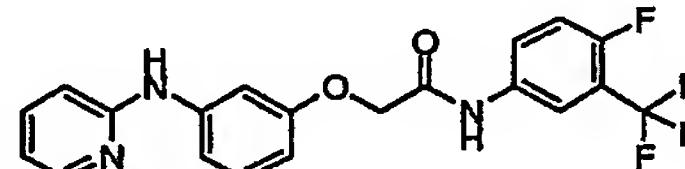
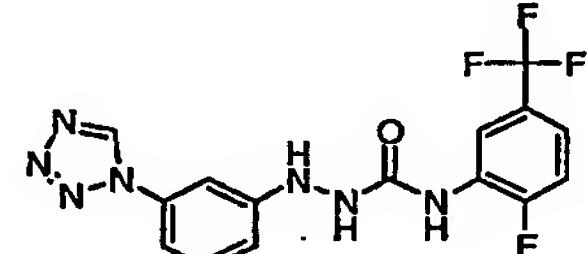
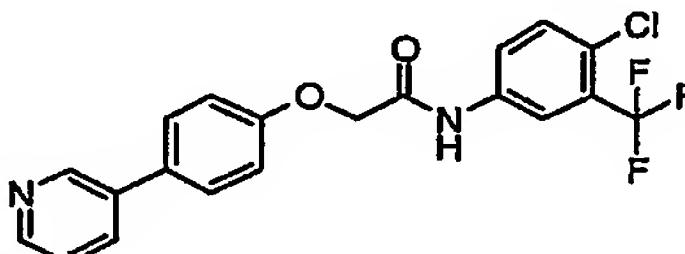
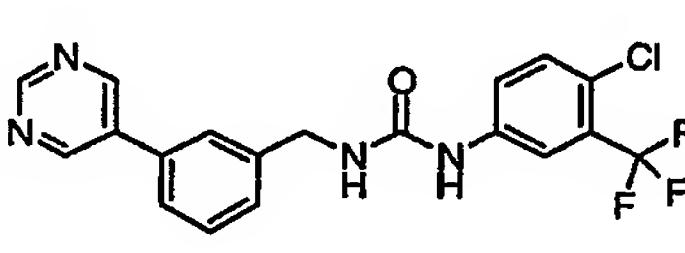
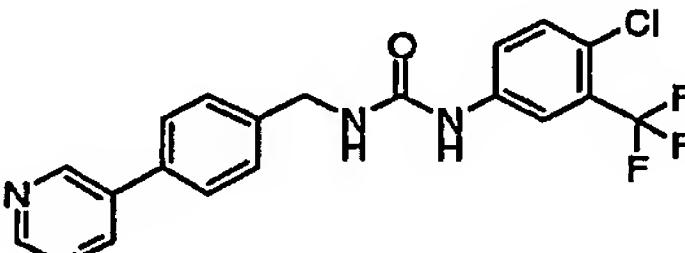
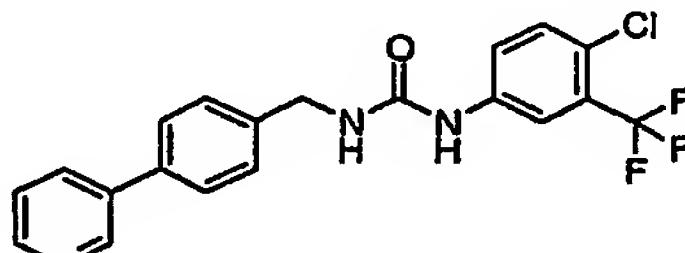
Entry	Name	Structure
77	N-[4-fluoro-3-(trifluoromethyl)phenyl]-N~2~-[3-(1H-tetrazol-1-yl)phenyl]glycinamide	
78	N-[4-fluoro-3-(trifluoromethyl)phenyl]-2-{[3-(pyridin-2-ylamino)phenyl]oxy}acetamide	
79	N-[2-fluoro-5-(trifluoromethyl)phenyl]-2-[3-(1H-tetrazol-1-yl)phenyl]hydrazinecarboxamide	
80	N-[4-chloro-3-(trifluoromethyl)phenyl]-2-[(4-pyridin-3-ylphenyl)oxy]acetamide	
81	N-[4-chloro-3-(trifluoromethyl)phenyl]-N'-(3-pyrimidin-5-ylphenyl)methylurea	
82	N-[4-chloro-3-(trifluoromethyl)phenyl]-N'-(4-pyrimidin-5-ylphenyl)methylurea	
83	N-[4-chloro-3-(trifluoromethyl)phenyl]-N'-(4-pyridin-3-ylphenyl)methylurea	

Table 3

Entry	Name	Structure
84	[3-(1H-tetrazol-1-yl)phenyl]methyl [4-chloro-3-(trifluoromethyl)phenyl]carbamate	
85	N-[4-fluoro-3-(trifluoromethyl)phenyl]-2-[(4-pyrimidin-5-ylphenyl)oxy]acetamide	
86	N~2~-[4-chloro-3-(trifluoromethyl)phenyl]-N-[3-(1H-tetrazol-1-yl)phenyl]glycinamide	
87	2-{[4-chloro-3-(trifluoromethyl)phenyl]oxy}-N-[3-(1H-tetrazol-1-yl)phenyl]acetamide	
88	N-[4-chloro-3-(trifluoromethyl)phenyl]-2-{{[3-methyl-4-(1H-tetrazol-1-yl)phenyl]oxy}acetamide}	
89	N-[4-chloro-3-(trifluoromethyl)phenyl]-2-{{[4-(1H-1,2,3-triazol-1-yl)phenyl]oxy}acetamide}	
90	N-[4-chloro-3-(trifluoromethyl)phenyl]-2-{{[3-fluoro-4-(1H-tetrazol-1-yl)phenyl]oxy}acetamide}	

Table 3

Entry	Name	Structure
91	N-[4-chloro-3-(trifluoromethyl)phenyl]-2-{[2-fluoro-4-(1H-tetrazol-1-yl)phenyl]oxy}acetamide	
92	N-({[4-chloro-3-(trifluoromethyl)phenyl]amino}carbonyl)-3-(1H-tetrazol-1-yl)benzenesulfonamide	
93	N-({[4-chloro-3-(trifluoromethyl)phenyl]amino}carbonyl)-N-methyl-3-(1H-tetrazol-1-yl)benzenesulfonamide	
94	N-[4-fluoro-3-(trifluoromethyl)phenyl]-2-[(4-pyridin-3-ylphenyl)oxy]acetamide	
95	2-({4-[2,4-bis(methoxy)pyrimidin-5-yl]phenyl}oxy)-N-[4-fluoro-3-(trifluoromethyl)phenyl]acetamide	
96	2-({4-[2,4-bis(methoxy)pyrimidin-5-yl]phenyl}oxy)-N-[4-chloro-3-(trifluoromethyl)phenyl]acetamide	
97	N-[4-chloro-3-(trifluoromethyl)phenyl]-2-[(4-pyridin-4-ylphenyl)oxy]acetamide	

Table 3

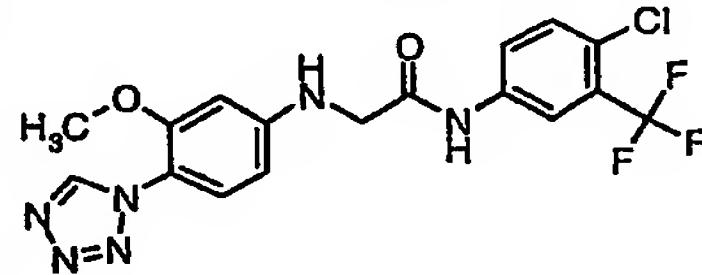
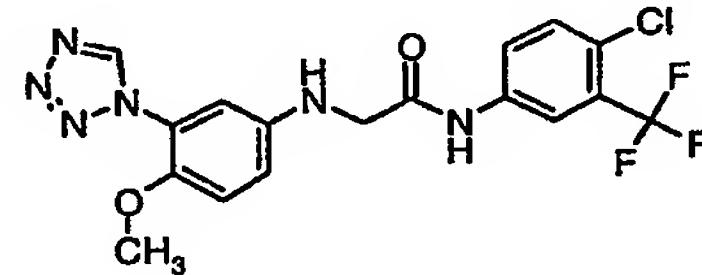
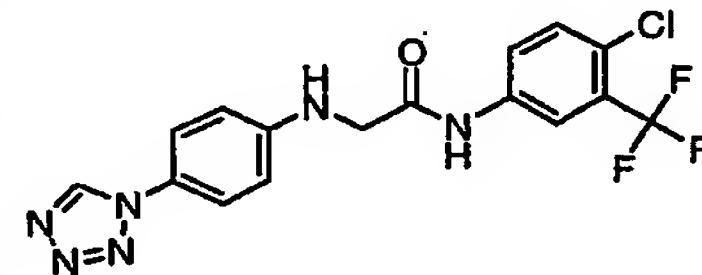
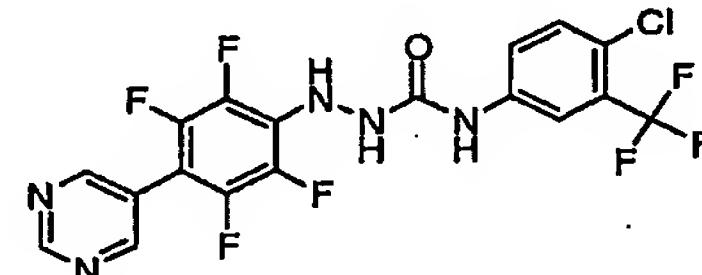
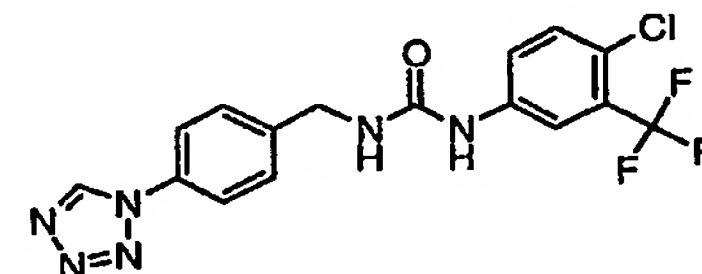
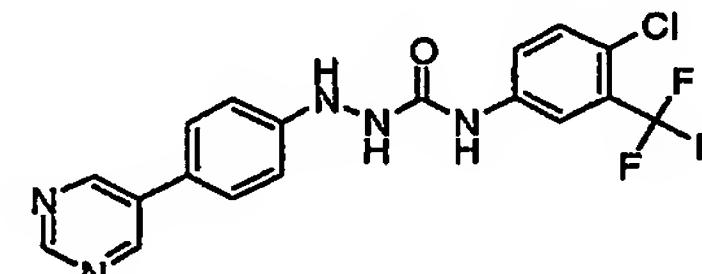
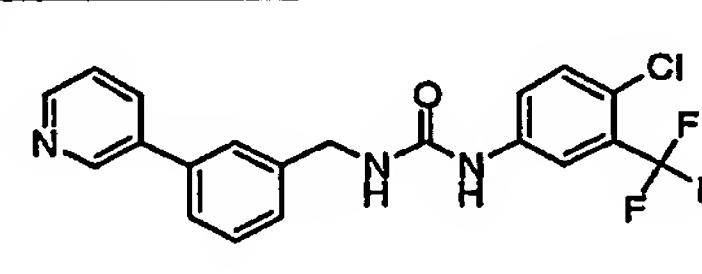
Entry	Name	Structure
98	N-[4-chloro-3-(trifluoromethyl)phenyl]-N~2~-[3-(methyloxy)-4-(1H-tetrazol-1-yl)phenyl]glycinamide	
99	N-[4-chloro-3-(trifluoromethyl)phenyl]-N~2~-[4-(methyloxy)-3-(1H-tetrazol-1-yl)phenyl]glycinamide	
100	N-[4-chloro-3-(trifluoromethyl)phenyl]-N~2~-[4-(1H-tetrazol-1-yl)phenyl]glycinamide	
101	N-[4-chloro-3-(trifluoromethyl)phenyl]-2-(2,3,5,6-tetrafluoro-4-pyrimidin-5-ylphenyl)hydrazinecarboxamide	
102	N-[4-chloro-3-(trifluoromethyl)phenyl]-N'-(4-(1H-tetrazol-1-yl)phenyl)methylurea	
103	N-[4-chloro-3-(trifluoromethyl)phenyl]-2-(4-pyrimidin-5-ylphenyl)hydrazinecarboxamide	
104	N-[4-chloro-3-(trifluoromethyl)phenyl]-N'-(3-pyridin-3-ylphenyl)methylurea	

Table 3

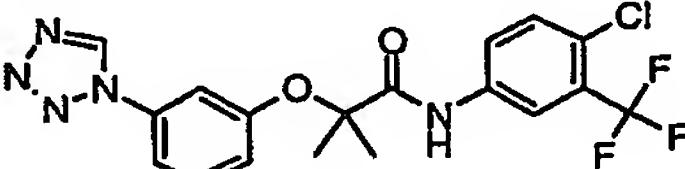
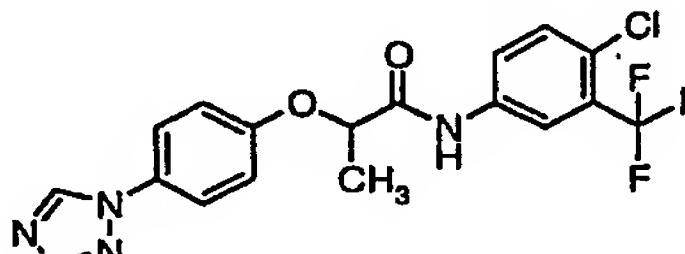
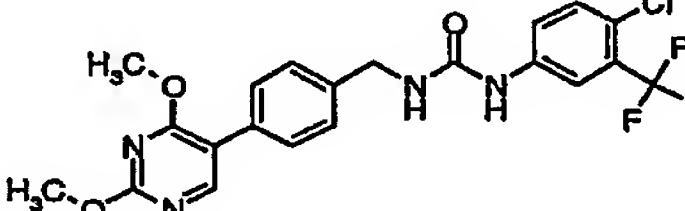
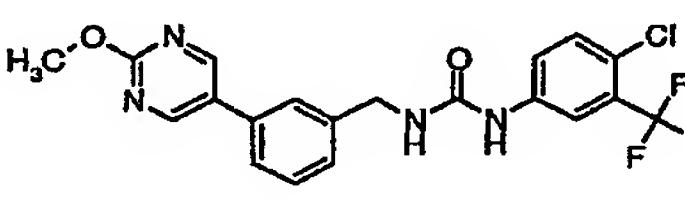
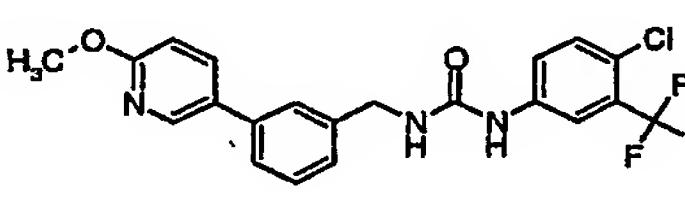
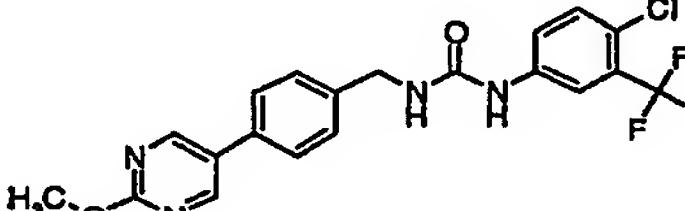
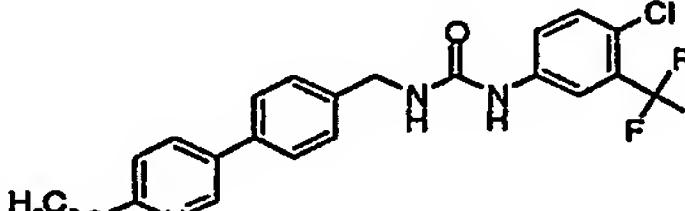
Entry	Name	Structure
105	N-[4-chloro-3-(trifluoromethyl)phenyl]-2-methyl-2-{[3-(1H-tetrazol-1-yl)phenyl]oxy}propanamide	
106	N-[4-chloro-3-(trifluoromethyl)phenyl]-2-{[4-(1H-tetrazol-1-yl)phenyl]oxy}propanamide	
107	N-({4-[2,4-bis(methyloxy)pyrimidin-5-yl]phenyl}methyl)-N'-[4-chloro-3-(trifluoromethyl)phenyl]urea	
108	N-[4-chloro-3-(trifluoromethyl)phenyl]-N'-({3-[2-(methyloxy)pyrimidin-5-yl]phenyl}methyl)urea	
109	N-[4-chloro-3-(trifluoromethyl)phenyl]-N'-({3-[6-(methyloxy)pyridin-3-yl]phenyl}methyl)urea	
110	N-[4-chloro-3-(trifluoromethyl)phenyl]-N'-({4-[2-(methyloxy)pyrimidin-5-yl]phenyl}methyl)urea	
111	N-[4-chloro-3-(trifluoromethyl)phenyl]-N'-({4-[6-(methyloxy)pyridin-3-yl]phenyl}methyl)urea	

Table 3

Entry	Name	Structure
112	1,1-dimethylethyl 2-{4-[2-{[4-chloro-3-(trifluoromethyl)phenyl]amino}-2-oxoethyl]oxy}phenyl}-1H-indole-1-carboxylate	
113	N-({[4-chloro-3-(trifluoromethyl)phenyl]amino}carbonyl)-4-(1H-tetrazol-1-yl)benzenesulfonamide	
114	N-[4-chloro-3-(trifluoromethyl)phenyl]-N~2~-[3-(2H-tetrazol-5-yl)phenyl]glycinamide	
115	N-[4-chloro-3-(trifluoromethyl)phenyl]-2-{{[2,6-difluoro-4-(1H-tetrazol-1-yl)phenyl]oxy}acetamide}	
116	(3-pyridin-3-ylphenyl)methyl [4-chloro-3-(trifluoromethyl)phenyl]carbamate	
117	(3-pyrimidin-5-ylphenyl)methyl [4-chloro-3-(trifluoromethyl)phenyl]carbamate	
118	(3-pyridin-4-ylphenyl)methyl [4-chloro-3-(trifluoromethyl)phenyl]carbamate	

Table 3

Entry	Name	Structure
119	N-[4-chloro-3-(trifluoromethyl)phenyl]-2-[4-(1H-tetrazol-1-yl)phenyl]hydrazinecarboxamide	
120	N-[4-chloro-3-(trifluoromethyl)phenyl]-2-(4-pyridin-3-ylphenyl)hydrazinecarboxamide	
121	(4-pyridin-3-ylphenyl)methyl [4-chloro-3-(trifluoromethyl)phenyl]carbamate	
122	(4-pyridin-4-ylphenyl)methyl [4-chloro-3-(trifluoromethyl)phenyl]carbamate	
123	(4-pyrimidin-5-ylphenyl)methyl [4-chloro-3-(trifluoromethyl)phenyl]carbamate	
124	N-[4-chloro-3-(trifluoromethyl)phenyl]-N'-(4-pyridin-4-ylphenyl)methylurea	
125	N-[4-chloro-3-(trifluoromethyl)phenyl]-2-(3-pyridin-3-ylphenyl)hydrazinecarboxamide	

Table 3

Entry	Name	Structure
126	N-[4-chloro-3-(trifluoromethyl)phenyl]-2-(3-pyrimidin-5-ylphenyl)hydrazinecarboxamide	
127	N-[5-chloro-2,4-bis(methoxy)phenyl]-N'-(4-pyrimidin-5-ylphenyl)methylurea	
128	N-[5-chloro-2,4-bis(methoxy)phenyl]-N'-(4-pyridin-3-ylphenyl)methylurea	
129	(4-pyrimidin-5-ylphenyl)methyl [5-chloro-2,4-bis(methoxy)phenyl]carbamate	
130	(4-pyridin-3-ylphenyl)methyl [5-chloro-2,4-bis(methoxy)phenyl]carbamate	
131	1-(4-pyridin-3-ylphenyl)ethyl [4-chloro-3-(trifluoromethyl)phenyl]carbamate	
132	1-(4-pyrimidin-5-ylphenyl)ethyl [4-chloro-3-(trifluoromethyl)phenyl]carbamate	

Table 3

Entry	Name	Structure
133	N-[5-chloro-2,4-bis(methyloxy)phenyl]-N'-(3-pyridin-3-ylphenyl)methyl]urea	
134	N-[5-chloro-2,4-bis(methyloxy)phenyl]-N'-(3-pyrimidin-5-ylphenyl)methyl]urea	
135	(3-pyridin-3-ylphenyl)methyl [5-chloro-2,4-bis(methyloxy)phenyl]carbamate	
136	(3-pyrimidin-5-ylphenyl)methyl [5-chloro-2,4-bis(methyloxy)phenyl]carbamate	
137	N-[4-chloro-3-(trifluoromethyl)phenyl]-2-methyl-2-(3-pyrimidin-5-ylphenyl)hydrazinecarboxamide	
138	N-[4-fluoro-3-(trifluoromethyl)phenyl]-N'-(4-pyridin-3-ylphenyl)methyl]urea	
139	N-{[3-(6-aminopyridin-3-yl)phenyl]methyl}-N'-(4-chloro-3-(trifluoromethyl)phenyl)urea	

Table 3

Entry	Name	Structure
140	N-{[4-(6-aminopyridin-3-yl)phenyl]methyl}-N'-[4-chloro-3-(trifluoromethyl)phenyl]urea	
141	N-{[3-(2-aminopyrimidin-5-yl)phenyl]methyl}-N'-[4-chloro-3-(trifluoromethyl)phenyl]urea	
142	N-{[4-(2-aminopyrimidin-5-yl)phenyl]methyl}-N'-[4-chloro-3-(trifluoromethyl)phenyl]urea	
143	N-[4-chloro-3-(trifluoromethyl)phenyl]-N'-[1-(4-pyridin-3-ylphenyl)ethyl]urea	
144	N-[4-chloro-3-(trifluoromethyl)phenyl]-N'-[1-(4-pyrimidin-5-ylphenyl)ethyl]urea	
145	N-[4-chloro-3-(trifluoromethyl)phenyl]-2-{{[4-(1H-indol-2-yl)phenyl]oxy}acetamide}	
146	N-[4-chloro-3-(trifluoromethyl)phenyl]-2-(isoquinolin-7-yloxy)acetamide	

Table 3

Entry	Name	Structure
147	N-[4-chloro-3-(trifluoromethyl)phenyl]-2-(4-pyridin-4-ylphenyl)hydrazinecarboxamide	
148	N-[4-chloro-3-(trifluoromethyl)phenyl]-2-(3-pyridin-4-ylphenyl)hydrazinecarboxamide	
149	N-[4-chloro-3-(trifluoromethyl)phenyl]-N'-(3-pyridin-4-ylphenyl)methylurea	
150	N-[4-chloro-3-(trifluoromethyl)phenyl]-N'-(3-quinoxalin-6-ylphenyl)methylurea	
151	methyl 3-amino-6-{[{[4-chloro-3-(trifluoromethyl)phenyl]amino}carbonyl]amino}methylphenylpyrazine-2-carboxylate	
152	N-[4-chloro-3-(trifluoromethyl)phenyl]-N'-(4-quinoxalin-6-ylphenyl)methylurea	
153	N-{[3-(2-amino-5-methylpyridin-3-yl)phenyl]methyl}-N'-[4-chloro-3-(trifluoromethyl)phenyl]urea	

Table 3

Entry	Name	Structure
154	methyl 3-amino-6-{{[[[4-chloro-3-(trifluoromethyl)phenyl]amino]carbonyl]amino]methyl}phenyl}pyrazine-2-carboxylate	
155	[3-(1H-tetrazol-1-yl)phenyl]methyl [3-chloro-4-(methyloxy)phenyl]carbamate	
156	N-[3-chloro-4-(methyloxy)phenyl]-N'-(3-(1H-tetrazol-1-yl)phenyl)methyl urea	
157	N-[4-chloro-3-(trifluoromethyl)phenyl]-2-{[4-(5-hydroxy-1H-tetrazol-1-yl)phenyl]oxy}acetamide	
158	N-{{[3-(2-amino-5-chloropyridin-3-yl)phenyl]methyl}-N'-[4-chloro-3-(trifluoromethyl)phenyl]urea	
159	N-{{[4-(2-amino-5-chloropyridin-3-yl)phenyl]methyl}-N'-[4-chloro-3-(trifluoromethyl)phenyl]urea	
160	N-{{[3-(6-chloropyridin-3-yl)phenyl]methyl}-N'-[4-chloro-3-(trifluoromethyl)phenyl]urea	

Table 3

Entry	Name	Structure
161	N-[4-(6-chloropyridin-3-yl)phenyl]methyl]-N'-[4-chloro-3-(trifluoromethyl)phenyl]urea	
162	N-[4-chloro-3-(trifluoromethyl)phenyl]-N'-{[4-(pyrimidin-2-yloxy)phenyl]methyl}urea	
163	N-({[4-chloro-3-(trifluoromethyl)phenyl]amino}carbonyl)-3-(1H-tetrazol-1-yl)benzamide	
164	3-amino-6-(3-{{[4-chloro-3-(trifluoromethyl)phenyl]amino}carbonyl}amino)methyl]phenyl)-N-[2-(dimethylamino)ethyl]pyrazine-2-carboxamide	
165	N-[4-chloro-3-(trifluoromethyl)phenyl]-N'-{[3-(6-fluoropyridin-3-yl)phenyl]methyl}urea	
166	N-[4-chloro-3-(trifluoromethyl)phenyl]-N'-{[3-[2-(methyloxy)pyridin-3-yl]phenyl]methyl}urea	
167	N-[4-chloro-3-(trifluoromethyl)phenyl]-N'-{[4-(6-fluoropyridin-3-yl)phenyl]methyl}urea	

Table 3

Entry	Name	Structure
168	N-[4-chloro-3-(trifluoromethyl)phenyl]-N'-(4-[2-(methyloxy)pyridin-3-yl]phenyl)methylurea	
169	N-[4-chloro-3-(trifluoromethyl)phenyl]-N'-(4-(6-methylpyridin-3-yl)phenyl)methylurea	
170	N-{[4-(2-amino-5-fluoropyridin-3-yl)phenyl]methyl}-N'-[4-chloro-3-(trifluoromethyl)phenyl]urea	
171	N-[4-chloro-3-(trifluoromethyl)phenyl]-N'-(3-(6-methylpyridin-3-yl)phenyl)methylurea	
172	N-{[4-(2-aminopyridin-3-yl)phenyl]methyl}-N'-[4-chloro-3-(trifluoromethyl)phenyl]urea	
173	N-{[3-(2-aminopyridin-3-yl)phenyl]methyl}-N'-[4-chloro-3-(trifluoromethyl)phenyl]urea	
174	[3-(6-methylpyridin-3-yl)phenyl]methyl [4-chloro-3-(trifluoromethyl)phenyl]carbamate	

Table 3

Entry	Name	Structure
175	[3-(2-amino-5-fluoropyridin-3-yl)phenyl]methyl [4-chloro-3-(trifluoromethyl)phenyl]carbamate	
176	[3-(2-aminopyridin-3-yl)phenyl]methyl [4-chloro-3-(trifluoromethyl)phenyl]carbamate	
177	(3-pyrazin-2-ylphenyl)methyl [4-chloro-3-(trifluoromethyl)phenyl]carbamate	
178	N-[4-chloro-3-(trifluoromethyl)phenyl]-N'-(3-[6-(hydroxymethyl)pyridin-3-yl]phenyl)methylurea	
179	N-[[3-(6-acetylpyridin-3-yl)phenyl]methyl]-N'-[4-chloro-3-(trifluoromethyl)phenyl]urea	
180	N-[4-chloro-3-(trifluoromethyl)phenyl]-N'-(3-(6-cyanopyridin-3-yl)phenyl)methylurea	
181	1,1-dimethylethyl (3S)-3-((3-amino-6-(3-((4-chloro-3-(trifluoromethyl)phenyl)amino)carbonyl)amino)methyl)phenyl)pyrazin-2-ylcarbonyl)piperidine-1-carboxylate	

Table 3

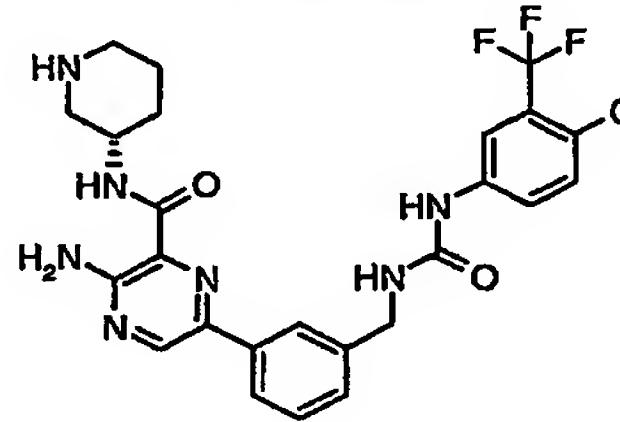
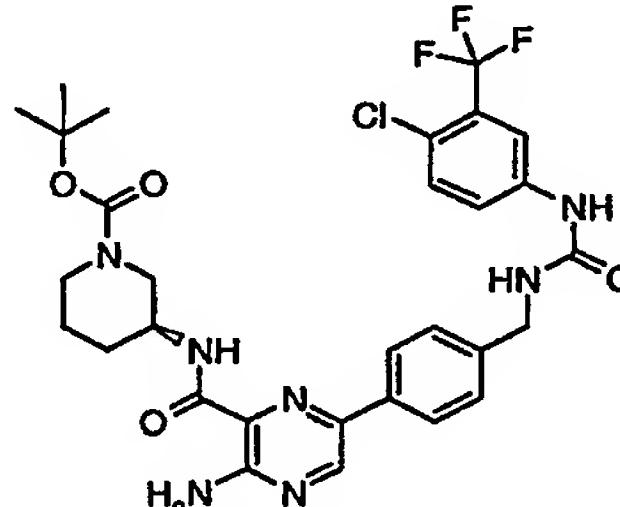
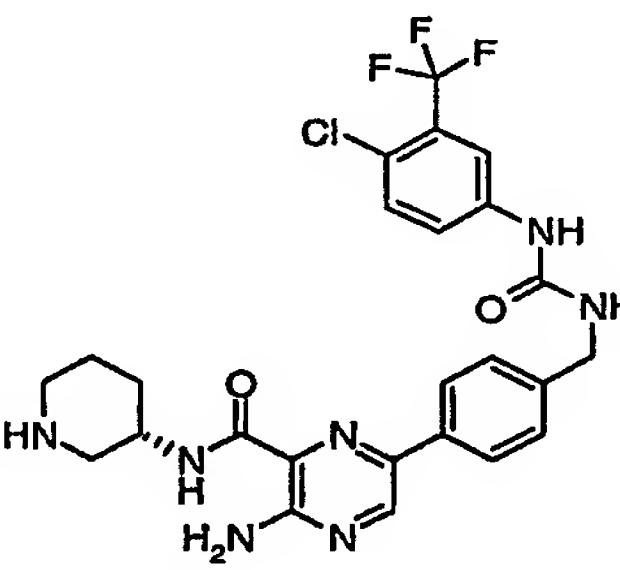
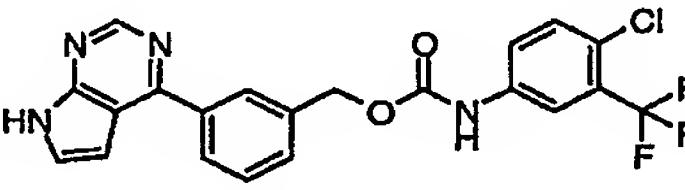
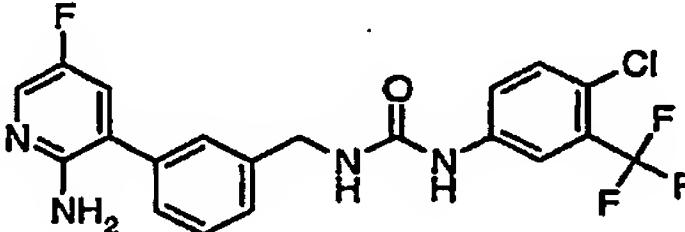
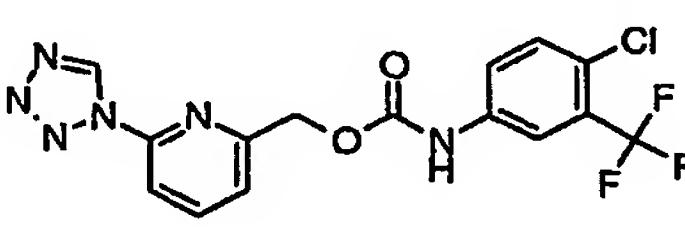
Entry	Name	Structure
182	3-amino-6-(3-{{[4-chloro-3-(trifluoromethyl)phenyl]amino}carbonyl}amino)methyl phenyl)-N-[ <i>(3S</i> )-piperidin-3-yl]pyrazine-2-carboxamide	
183	1,1-dimethylethyl ( <i>3S</i> )-3-({[3-amino-6-(4-((4-chloro-3-(trifluoromethyl)phenyl)amino)carbonyl)amino]methyl}phenyl)pyrazin-2-yl]carbonyl}amino)piperidine-1-carboxylate	
184	3-amino-6-(4-{{[4-chloro-3-(trifluoromethyl)phenyl]amino}carbonyl}amino)methyl phenyl)-N-[ <i>(3S</i> )-piperidin-3-yl]pyrazine-2-carboxamide	
185	[3-(7 <i>H</i> -pyrrolo[2,3- <i>d</i> ]pyrimidin-4-yl)phenyl]methyl [4-chloro-3-(trifluoromethyl)phenyl]carbamate	
186	N-{{[3-(2-amino-5-fluoropyridin-3-yl)phenyl]methyl}-N'-[4-chloro-3-(trifluoromethyl)phenyl]urea	
187	[6-(1 <i>H</i> -tetrazol-1-yl)pyridin-2-yl]methyl [4-chloro-3-(trifluoromethyl)phenyl]carbamate	

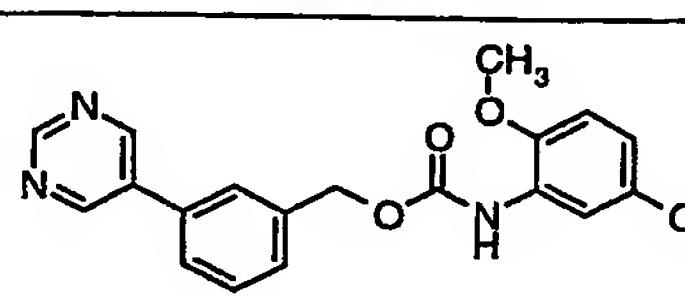
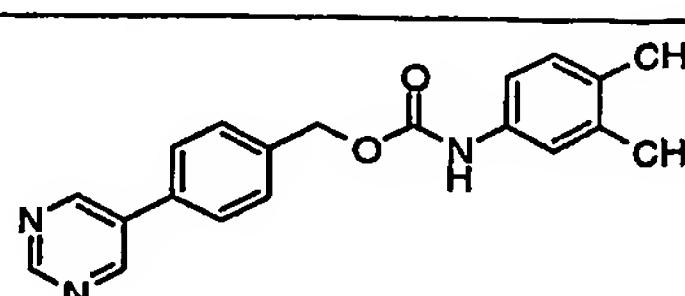
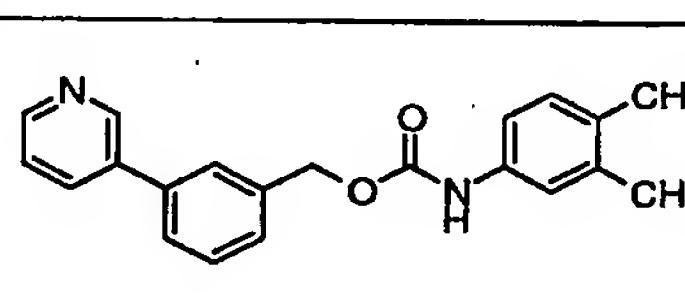
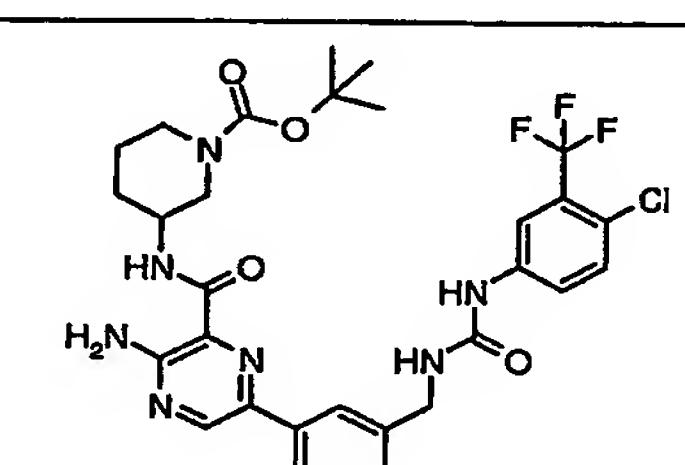
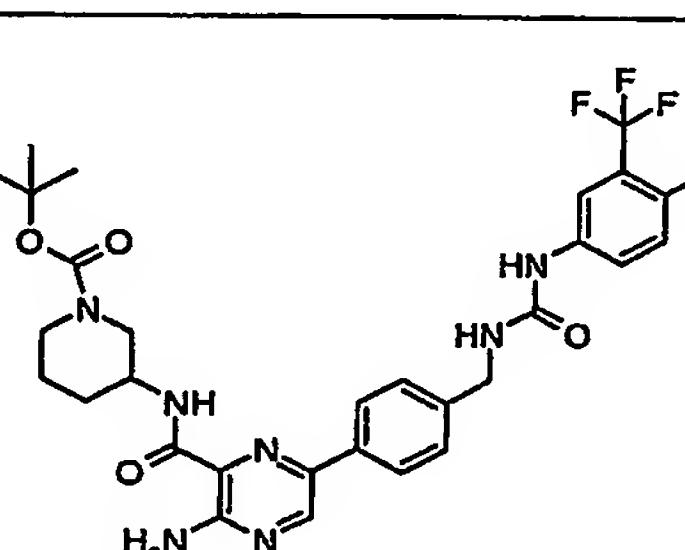
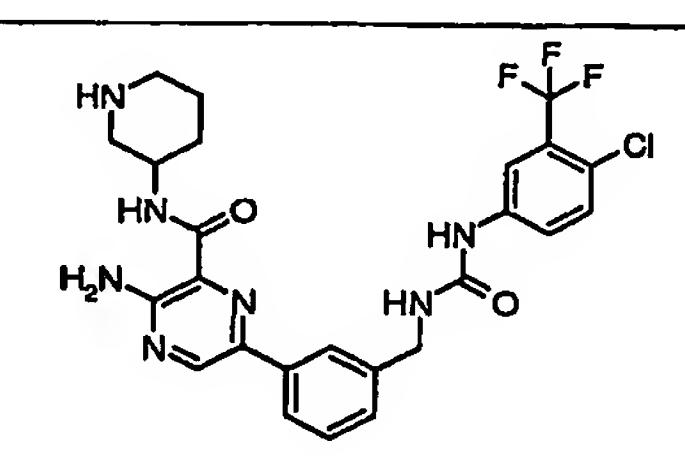
Table 3

Entry	Name	Structure
188	[3-(1H-benzimidazol-2-yl)phenyl]methyl [4-chloro-3-(trifluoromethyl)phenyl]carbamate	
189	[3-(6-amino-2-methylpyridin-3-yl)phenyl]methyl [4-chloro-3-(trifluoromethyl)phenyl]carbamate	
190	N-[4-chloro-3-(trifluoromethyl)phenyl]-N'-(3-[5-(methylthio)pyridin-3-yl]phenyl)methylurea	
191	[4-(6-methylpyridin-3-yl)phenyl]methyl [4-chloro-3-(trifluoromethyl)phenyl]carbamate	
192	[4-(2-amino-5-fluoropyridin-3-yl)phenyl]methyl [4-chloro-3-(trifluoromethyl)phenyl]carbamate	
193	[4-(2-aminopyridin-3-yl)phenyl]methyl [4-chloro-3-(trifluoromethyl)phenyl]carbamate	
194	(4-pyrazin-2-ylphenyl)methyl [4-chloro-3-(trifluoromethyl)phenyl]carbamate	

Table 3

Entry	Name	Structure
195	[4-(7H-pyrrolo[2,3-d]pyrimidin-4-yl)phenyl]methyl [4-chloro-3-(trifluoromethyl)phenyl]carbamate	
196	[4-(6-amino-2-methylpyridin-3-yl)phenyl]methyl [4-chloro-3-(trifluoromethyl)phenyl]carbamate	
197	[3-(1H-tetrazol-1-yl)phenyl]methyl 1,3-benzothiazol-2-ylcarbamate	
198	[3-(1H-tetrazol-1-yl)phenyl]methyl (5-bromopyridin-2-yl)carbamate	
199	(3-pyridin-3-ylphenyl)methyl (3,5-dimethylphenyl)carbamate	
200	(3-pyridin-3-ylphenyl)methyl [5-chloro-2-(methyoxy)phenyl]carbamate	
201	[4-(1H-tetrazol-1-yl)phenyl]methyl [4-chloro-3-(trifluoromethyl)phenyl]carbamate	

Table 3

Entry	Name	Structure
202	(3-pyrimidin-5-ylphenyl)methyl [5-chloro-2-(methyoxy)phenyl]carbamate	
203	(4-pyrimidin-5-ylphenyl)methyl (3,4-dimethylphenyl)carbamate	
204	(3-pyridin-3-ylphenyl)methyl (3,4-dimethylphenyl)carbamate	
205	1,1-dimethylethyl 3-({[3-amino-6-(3-[(4-chloro-3-(trifluoromethyl)phenyl]amino)carbonyl]amino]methyl}phenyl)pyrazin-2-yl]carbonyl}amino)piperidine-1-carboxylate	
206	1,1-dimethylethyl 3-({[3-amino-6-(4-[(4-chloro-3-(trifluoromethyl)phenyl]amino)carbonyl]amino]methyl}phenyl)pyrazin-2-yl]carbonyl}amino)piperidine-1-carboxylate	
207	3-amino-6-(3-{[(4-chloro-3-(trifluoromethyl)phenyl]amino}carbonyl)amino]methyl}phenyl)-N-piperidin-3-ylpyrazine-2-carboxamide	

**Table 3**

Entry	Name	Structure
208	3-amino-6-{[({{4-chloro-3-(trifluoromethyl)phenyl}amino}carbonyl)amino]methyl}phenyl)-N-piperidin-3-ylpyrazine-2-carboxamide	
209	1,1-dimethylethyl 4-{[3-amino-6-(3-(({{4-chloro-3-(trifluoromethyl)phenyl}amino}carbonyl)amino)methyl)phenyl]pyrazin-2-yl]carbonyl}piperazine-1-carboxylate	
210	1,1-dimethylethyl 4-{[3-amino-6-(4-(({{4-chloro-3-(trifluoromethyl)phenyl}amino}carbonyl)amino)methyl)phenyl]pyrazin-2-yl]carbonyl}piperazine-1-carboxylate	
211	N-(3-[5-amino-6-(piperazin-1-ylcarbonyl)pyrazin-2-yl]phenyl)methyl-N'-[4-chloro-3-(trifluoromethyl)phenyl]urea	

Table 3

Entry	Name	Structure
212	N-({4-[5-amino-6-(piperazin-1-ylcarbonyl)pyrazin-2-yl]phenyl}methyl)-N'-[4-chloro-3-(trifluoromethyl)phenyl]urea	
213	N-[4-chloro-3-(trifluoromethyl)phenyl]-N'-{[3-(1H-pyrazol-4-yl)phenyl]methyl}urea	
214	N-[4-chloro-3-(trifluoromethyl)phenyl]-N'-{[4-(1H-pyrazol-4-yl)phenyl]methyl}urea	
215	[3-(2-piperazin-1-ylpyrimidin-5-yl)phenyl]methyl [4-chloro-3-(trifluoromethyl)phenyl]carbamate	
216	[4-(2-piperazin-1-ylpyrimidin-5-yl)phenyl]methyl [4-chloro-3-(trifluoromethyl)phenyl]carbamate	
217	N-{{3-(2-chloropyridin-3-yl)phenyl}methyl}-N'-[4-chloro-3-(trifluoromethyl)phenyl]urea	

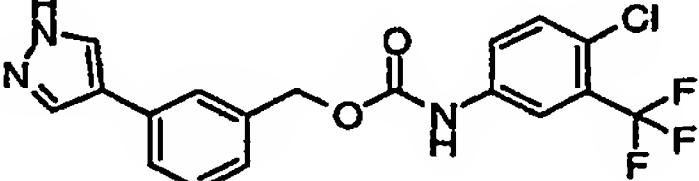
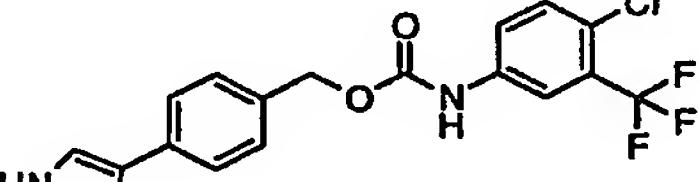
Table 3

Entry	Name	Structure
218	N-{[4-(2-chloropyridin-3-yl)phenyl]methyl}-N'-[4-chloro-3-(trifluoromethyl)phenyl]urea	
219	N-[4-chloro-3-(trifluoromethyl)phenyl]-N'-{[3-(2-fluoropyridin-3-yl)phenyl]methyl}urea	
220	N-[4-chloro-3-(trifluoromethyl)phenyl]-N'-{[4-(2-fluoropyridin-3-yl)phenyl]methyl}urea	
221	[3-(1H-tetrazol-1-yl)phenyl]methyl [3-(trifluoromethyl)phenyl]carbamate	
222	[3-(1H-tetrazol-1-yl)phenyl]methyl [6-(trifluoromethyl)pyridin-2-yl]carbamate	
223	[3-(1H-tetrazol-1-yl)phenyl]methyl [4-(trifluoromethyl)pyridin-2-yl]carbamate	
224	N-[4-chloro-3-(trifluoromethyl)phenyl]-N'-{[3-[5-(methylthio)pyridin-2-yl]phenyl]methyl}urea	

Table 3

Entry	Name	Structure
225	[3-(2,6-dimethylpyridin-3-yl)phenyl]methyl [4-chloro-3-(trifluoromethyl)phenyl]carbamate	
226	{3-[5-(methyloxy)pyridin-3-yl]phenyl}methyl [4-chloro-3-(trifluoromethyl)phenyl]carbamate	
227	2,3'-bipyridin-6-ylmethyl [4-chloro-3-(trifluoromethyl)phenyl]carbamate	
228	(6-pyrimidin-5-ylpyridin-2-yl)methyl [4-chloro-3-(trifluoromethyl)phenyl]carbamate	
229	N-[4-chloro-3-(trifluoromethyl)phenyl]-N'-(3-isoquinolin-4-ylphenyl)methylurea	
230	N-[4-chloro-3-(trifluoromethyl)phenyl]-N'-(4-isoquinolin-4-ylphenyl)methylurea	
231	[6-(1H-tetrazol-1-yl)pyridin-2-yl]methyl [4-(trifluoromethyl)pyridin-2-yl]carbamate	

Table 3

Entry	Name	Structure
232	[3-(1H-pyrazol-4-yl)phenyl]methyl [4-chloro-3-(trifluoromethyl)phenyl]carbamate	
233	[4-(1H-pyrazol-4-yl)phenyl]methyl [4-chloro-3-(trifluoromethyl)phenyl]carbamate	

28. A pharmaceutical composition comprising the compound according to any one of claims 1 - 27 and a pharmaceutically acceptable carrier.

29. A metabolite of the compound or the pharmaceutical composition according to any one of claims 1 - 28.

30. A method for modulating the *in-vivo* activity of a kinase, the method comprising administering to a subject an effective amount of the compound according to any of claims 1 - 27 or a compound selected from N-naphthalen-1-yl-2-{[3-(1H-tetrazol-1-yl)phenyl]oxy} acetamide, N-[4-(phenyloxy)phenyl]-2-{[3-(1H-tetrazol-1-yl)phenyl]oxy} acetamide, N-(3,4-dimethylphenyl)-2-{[3-(1H-tetrazol-1-yl)phenyl]oxy} acetamide, N-(2,3-dimethylphenyl)-2-{[3-(1H-tetrazol-1-yl)phenyl]oxy} acetamide, N-(2,4-dimethylphenyl)-2-{[3-(1H-tetrazol-1-yl)phenyl]oxy} acetamide, N-(2,5-dimethylphenyl)-2-{[3-(1H-tetrazol-1-yl)phenyl]oxy}acetamide, N-(3,5-dimethylphenyl)-2-{[3-(1H-tetrazol-1-yl)phenyl]oxy} acetamide, N-(2,6-dimethyl-phenyl)-2-{[3-(1H-tetrazol-1-yl)phenyl]oxy}acetamide, 2-{[3-(1H-tetrazol-1-yl)phenyl]oxy}-N-(2,4,6-trimethylphenyl)acetamide, N-(2-ethylphenyl)-2-{[3-(1H-tetrazol-1-yl)phenyl]oxy} acetamide, N-(4-ethylphenyl)-2-{[3-(1H-tetrazol-1-yl)phenyl]oxy} acetamide, N-(2,6-diethylphenyl)-2-{[3-(1H-tetrazol-1-yl)phenyl]oxy}acetamide, N-[2-(methyloxy)phenyl]-2-{[3-(1H-tetrazol-1-yl)phenyl]oxy} acetamide, N-[2-(ethyloxy) phenyl]-2-{[3-(1H-tetrazol-1-yl)phenyl]oxy}acetamide, N-[3-(ethyloxy)phenyl]-2-{[3-(1H-tetrazol-1-yl)phenyl]oxy} acetamide, N-[2,4-bis(methyloxy)phenyl]-2-{[3-(1H-tetrazol-1-

yl)phenyl]oxy} acetamide, N-[4-(dimethylamino)phenyl]-2-{[3-(1H-tetrazol-1-yl)phenyl]oxy}acetamide, N-(2,3-dichlorophenyl)-2-{[3-(1H-tetrazol-1-yl)phenyl]oxy}acetamide, N-(4-chloro-3-methylphenyl)-2-{[3-(1H-tetrazol-1-yl)phenyl]oxy}acetamide, N-(4-bromophenyl)-2-{[3-(1H-tetrazol-1-yl)phenyl]oxy}acetamide, N-(2-fluorophenyl)-2-{[3-(1H-tetrazol-1-yl)phenyl]oxy}acetamide, N-(4-fluorophenyl)-2-{[3-(1H-tetrazol-1-yl)phenyl]oxy}acetamide, 2-{[3-(1H-tetrazol-1-yl)phenyl]oxy}-N-[2-(trifluoromethyl)phenyl]acetamide, 2-{[3-(1H-tetrazol-1-yl)phenyl]oxy}-N-[3-(trifluoromethyl)phenyl]acetamide, methyl 4-[({[3-(1H-tetrazol-1-yl)phenyl]oxy}acetyl)amino] benzoate, ethyl 4-[({[3-(1H-tetrazol-1-yl)phenyl]oxy}acetyl)amino] benzoate, 3-[({[3-(1H-tetrazol-1-yl)phenyl]oxy}acetyl)amino] benzoic acid, N-[3-(methyloxy)phenyl]-2-{[3-(1H-tetrazol-1-yl)phenyl]oxy}acetamide, N-[4-(methyloxy)phenyl]-2-{[3-(1H-tetrazol-1-yl)phenyl]oxy}acetamide, N-[2-chloro-5-(trifluoromethyl)phenyl]-2-{[3-(1H-tetrazol-1-yl)phenyl]oxy}acetamide, N-[4-chloro-3-(trifluoromethyl)phenyl]-2-{[3-(4H-1,2,4-triazol-4-yl)phenyl]oxy}acetamide, N-(4-chlorophenyl)-2-{[3-(1H-tetrazol-1-yl)phenyl]oxy}acetamide, N-(4-aminophenyl)-2-{[3-(1H-tetrazol-1-yl)phenyl]oxy}acetamide, and N-(4-acetylphenyl)-2-{[3-(1H-tetrazol-1-yl)phenyl]oxy}acetamide.

31. The method according to claim 30, wherein the kinase is c-Kit.
32. The method according to claim 31, wherein modulating the *in vivo* activity of c-Kit comprises inhibition of c-Kit.
33. A method of treating diseases or disorders associated with uncontrolled, abnormal, and/or unwanted cellular activities, the method comprising administering, to a mammal in need thereof, a therapeutically effective amount of the compound or the pharmaceutical composition as described in any one of claims 1 -28 or a compound, or a pharmaceutical composition comprising said compound, selected from N-naphthalen-1-yl-2-{[3-(1H-tetrazol-1-yl)phenyl]oxy}acetamide, N-[4-(phenyloxy)phenyl]-2-{[3-(1H-tetrazol-1-yl)phenyl]oxy}acetamide, N-(3,4-dimethylphenyl)-2-{[3-(1H-tetrazol-1-yl)phenyl]oxy}acetamide, N-(2,3-dimethyl-phenyl)-2-{[3-(1H-tetrazol-1-yl)phenyl]oxy}acetamide, N-(2,4-dimethylphenyl)-2-{[3-(1H-tetrazol-1-yl)phenyl]oxy}acetamide, N-(2,5-dimethylphenyl)-2-{[3-(1H-tetrazol-1-yl)phenyl]oxy}acetamide, N-(3,5-dimethylphenyl)-2-{[3-(1H-tetrazol-1-yl)phenyl]oxy}acetamide, N-(2,6-dimethyl-phenyl)-2-{[3-(1H-tetrazol-1-yl)phenyl]oxy}acetamide, 2-{[3-(1H-tetrazol-1-yl)phenyl]oxy}-N-(2,4,6-trimethylphenyl)acetamide, N-(2-ethyl-phenyl)-2-{[3-(1H-tetrazol-1-yl)phenyl]oxy}

acetamide, N-(4-ethylphenyl)-2-{[3-(1H-tetrazol-1-yl)phenyl]oxy} acetamide, N-(2,6-diethylphenyl)-2-{[3-(1H-tetrazol-1-yl)phenyl]oxy} acetamide, N-[2-(methyloxy)phenyl]-2-{[3-(1H-tetrazol-1-yl)phenyl]oxy} acetamide, N-[2-(ethyloxy) phenyl]-2-{[3-(1H-tetrazol-1-yl)phenyl]oxy}acetamide, N-[3-(ethyloxy)phenyl]-2-{[3-(1H-tetrazol-1-yl)phenyl]oxy} acetamide, N-[2,4-bis(methyl-oxy)phenyl]-2-{[3-(1H-tetrazol-1-yl)phenyl]oxy} acetamide, N-[4-(dimethylamino)-phenyl]-2-{[3-(1H-tetrazol-1-yl)phenyl]oxy}acetamide, N-(2,3-dichlorophenyl)-2-{[3-(1H-tetrazol-1-yl)phenyl]oxy} acetamide, N-(4-chloro-3-methylphenyl)-2-{[3-(1H-tetrazol-1-yl)phenyl]oxy} acetamide, N-(4-bromophenyl)-2-{[3-(1H-tetrazol-1-yl)phenyl]oxy} acetamide, N-(2-fluorophenyl)-2-{[3-(1H-tetrazol-1-yl)phenyl]oxy} acetamide, N-(4-fluorophenyl)-2-{[3-(1H-tetrazol-1-yl)phenyl]oxy} acetamide, 2-{[3-(1H-tetrazol-1-yl)phenyl]oxy}-N-[2-(trifluoromethyl)phenyl] acetamide, 2-{[3-(1H-tetrazol-1-yl)phenyl]oxy}-N-[3-(trifluoromethyl)phenyl] acetamide, methyl 4-[( {[3-(1H-tetrazol-1-yl)phenyl]oxy} acetyl)amino] benzoate, ethyl 4-[( {[3-(1H-tetrazol-1-yl)phenyl]oxy} acetyl)amino] benzoate, 3-[( {[3-(1H-tetrazol-1-yl)phenyl]oxy} acetyl) amino] benzoic acid, N-[3-(methyloxy)phenyl]-2-{[3-(1H-tetrazol-1-yl)phenyl]oxy} acetamide, N-[4-(methyloxy)phenyl]-2-{[3-(1H-tetrazol-1-yl)phenyl]oxy} acetamide, N-[2-chloro-5-(trifluoromethyl)phenyl]-2-{[3-(1H-tetrazol-1-yl)phenyl]oxy} acetamide, N-[4-chloro-3-(trifluoromethyl)phenyl]-2-{[3-(4H-1,2,4-triazol-4-yl)phenyl]oxy} acetamide, N-(4-chlorophenyl)-2-{[3-(1H-tetrazol-1-yl)phenyl]oxy} acetamide, N-(4-aminophenyl)-2-{[3-(1H-tetrazol-1-yl)phenyl]oxy} acetamide, and N-(4-acetylphenyl)-2-{[3-(1H-tetrazol-1-yl)phenyl]oxy} acetamide.

34. A method of screening for modulators of c-Kit, the method comprising combining the compound according to any one of claims 1 -27 or a compound selected from N-naphthalen-1-yl-2-{[3-(1H-tetrazol-1-yl)phenyl]oxy} acetamide, N-[4-(phenyloxy)phenyl]-2-{[3-(1H-tetrazol-1-yl)phenyl]oxy} acetamide, N-(3,4-dimethylphenyl)-2-{[3-(1H-tetrazol-1-yl)phenyl]oxy} acetamide, N-(2,3-dimethylphenyl)-2-{[3-(1H-tetrazol-1-yl)phenyl]oxy} acetamide, N-(2,4-dimethylphenyl)-2-{[3-(1H-tetrazol-1-yl)phenyl]oxy} acetamide, N-(2,5-dimethylphenyl)-2-{[3-(1H-tetrazol-1-yl)phenyl]oxy} acetamide, N-(3,5-dimethylphenyl)-2-{[3-(1H-tetrazol-1-yl)phenyl]oxy} acetamide, N-(2,6-dimethylphenyl)-2-{[3-(1H-tetrazol-1-yl)phenyl]oxy}acetamide, 2-{[3-(1H-tetrazol-1-yl)phenyl]oxy}-N-(2,4,6-trimethylphenyl)acetamide, N-(2-ethylphenyl)-2-{[3-(1H-tetrazol-1-yl)phenyl]oxy} acetamide, N-(4-ethylphenyl)-2-{[3-(1H-tetrazol-1-yl)phenyl]oxy} acetamide, N-(2,6-diethylphenyl)-2-{[3-(1H-tetrazol-1-yl)phenyl]oxy}acetamide, N-[2-

(methyloxy)phenyl]-2-{[3-(1H-tetrazol-1-yl)phenyl]oxy} acetamide, N-[2-(ethyloxy)phenyl]-2-{[3-(1H-tetrazol-1-yl)phenyl]oxy}acetamide, N-[3-(ethyloxy)phenyl]-2-{[3-(1H-tetrazol-1-yl)phenyl]oxy} acetamide, N-[2,4-bis(methyloxy)phenyl]-2-{[3-(1H-tetrazol-1-yl)phenyl]oxy} acetamide, N-[4-(dimethylamino)phenyl]-2-{[3-(1H-tetrazol-1-yl)phenyl]oxy}acetamide, N-(2,3-dichlorophenyl)-2-{[3-(1H-tetrazol-1-yl)phenyl]oxy}acetamide, N-(4-chloro-3-methylphenyl)-2-{[3-(1H-tetrazol-1-yl)phenyl]oxy}acetamide, N-(4-bromophenyl)-2-{[3-(1H-tetrazol-1-yl)phenyl]oxy}acetamide, N-(2-fluorophenyl)-2-{[3-(1H-tetrazol-1-yl)phenyl]oxy} acetamide, N-(4-fluorophenyl)-2-{[3-(1H-tetrazol-1-yl)phenyl]oxy}acetamide, 2-{[3-(1H-tetrazol-1-yl)phenyl]oxy}-N-[2-(trifluoromethyl)phenyl] acetamide, 2-{[3-(1H-tetrazol-1-yl)phenyl]oxy}-N-[3-(trifluoromethyl)phenyl] acetamide, methyl 4-[({[3-(1H-tetrazol-1-yl)phenyl]oxy}acetyl)amino] benzoate, ethyl 4-[({[3-(1H-tetrazol-1-yl)phenyl]oxy}acetyl)amino] benzoate, 3-[({[3-(1H-tetrazol-1-yl)phenyl]oxy}acetyl)amino] benzoic acid, N-[3-(methyloxy)phenyl]-2-{[3-(1H-tetrazol-1-yl)phenyl]oxy} acetamide, N-[4-(methyloxy)phenyl]-2-{[3-(1H-tetrazol-1-yl)phenyl]oxy} acetamide, N-[2-chloro-5-(trifluoromethyl)phenyl]-2-{[3-(1H-tetrazol-1-yl)phenyl]oxy} acetamide, N-[4-chloro-3-(trifluoromethyl)phenyl]-2-{[3-(4H-1,2,4-triazol-4-yl)phenyl]oxy} acetamide, N-(4-chlorophenyl)-2-{[3-(1H-tetrazol-1-yl)phenyl]oxy} acetamide, N-(4-aminophenyl)-2-{[3-(1H-tetrazol-1-yl)phenyl]oxy} acetamide, and N-(4-acetylphenyl)-2-{[3-(1H-tetrazol-1-yl)phenyl]oxy} acetamide, and at least one candidate agent and determining the effect of the candidate agent on c-Kit activity.

35. A method of inhibiting proliferative activity in a cell, the method comprising administering an effective amount of a composition comprising the compound according any one of claims 1 - 27 or a compound selected from N-naphthalen-1-yl-2-{[3-(1H-tetrazol-1-yl)phenyl]oxy} acetamide, N-[4-(phenyloxy)phenyl]-2-{[3-(1H-tetrazol-1-yl)phenyl]oxy} acetamide, N-(3,4-dimethylphenyl)-2-{[3-(1H-tetrazol-1-yl)phenyl]oxy} acetamide, N-(2,3-dimethylphenyl)-2-{[3-(1H-tetrazol-1-yl)phenyl]oxy} acetamide, N-(2,4-dimethylphenyl)-2-{[3-(1H-tetrazol-1-yl)phenyl]oxy} acetamide, N-(2,5-dimethylphenyl)-2-{[3-(1H-tetrazol-1-yl)phenyl]oxy}acetamide, N-(3,5-dimethylphenyl)-2-{[3-(1H-tetrazol-1-yl)phenyl]oxy} acetamide, N-(2,6-dimethyl-phenyl)-2-{[3-(1H-tetrazol-1-yl)phenyl]oxy}acetamide, N-(2,6-dimethylphenyl)-2-{[3-(1H-tetrazol-1-yl)phenyl]oxy}acetamide, 2-{[3-(1H-tetrazol-1-yl)phenyl]oxy}-N-(2,4,6-trimethylphenyl)acetamide, N-(2-ethylphenyl)-2-{[3-(1H-tetrazol-1-yl)phenyl]oxy} acetamide, N-(4-ethylphenyl)-2-{[3-(1H-tetrazol-1-yl)phenyl]oxy} acetamide, N-(2,6-diethylphenyl)-2-{[3-(1H-tetrazol-1-yl)phenyl]oxy}acetamide, N-[2-(methyloxy)phenyl]-

2-{{[3-(1H-tetrazol-1-yl)phenyl]oxy} acetamide, N-[2-(ethyloxy) phenyl]-2-{{[3-(1H-tetrazol-1-yl)phenyl]oxy}acetamide, N-[3-(ethyloxy)phenyl]-2-{{[3-(1H-tetrazol-1-yl)phenyl]oxy} acetamide, N-[2,4-bis(methyloxy)phenyl]-2-{{[3-(1H-tetrazol-1-yl)phenyl]oxy} acetamide, N-[4-(dimethylamino)phenyl]-2-{{[3-(1H-tetrazol-1-yl)phenyl]oxy}acetamide, N-(2,3-dichlorophenyl)-2-{{[3-(1H-tetrazol-1-yl)phenyl]oxy}acetamide, N-(4-chloro-3-methylphenyl)-2-{{[3-(1H-tetrazol-1-yl)phenyl]oxy}acetamide, N-(4-bromophenyl)-2-{{[3-(1H-tetrazol-1-yl)phenyl]oxy}acetamide, N-(2-fluorophenyl)-2-{{[3-(1H-tetrazol-1-yl)phenyl]oxy} acetamide, N-(4-fluorophenyl)-2-{{[3-(1H-tetrazol-1-yl)phenyl]oxy} acetamide, 2-{{[3-(1H-tetrazol-1-yl)phenyl]oxy}-N-[2-(trifluoromethyl)phenyl] acetamide, 2-{{[3-(1H-tetrazol-1-yl)phenyl]oxy}-N-[3-(trifluoromethyl)phenyl] acetamide, methyl 4-[({{[3-(1H-tetrazol-1-yl)phenyl]oxy}acetyl)amino] benzoate, ethyl 4-[({{[3-(1H-tetrazol-1-yl)phenyl]oxy}acetyl)amino] benzoate, 3-[({{[3-(1H-tetrazol-1-yl)phenyl]oxy}acetyl)amino] benzoic acid, N-[3-(methyloxy)phenyl]-2-{{[3-(1H-tetrazol-1-yl)phenyl]oxy} acetamide, N-[4-(methyloxy)phenyl]-2-{{[3-(1H-tetrazol-1-yl)phenyl]oxy} acetamide, N-[2-chloro-5-(trifluoromethyl)phenyl]-2-{{[3-(1H-tetrazol-1-yl)phenyl]oxy} acetamide, N-[4-chloro-3-(trifluoromethyl)phenyl]-2-{{[3-(4H-1,2,4-triazol-4-yl)phenyl]oxy} acetamide, N-(4-chlorophenyl)-2-{{[3-(1H-tetrazol-1-yl)phenyl]oxy} acetamide, N-(4-aminophenyl)-2-{{[3-(1H-tetrazol-1-yl)phenyl]oxy} acetamide, and N-(4-acetylphenyl)-2-{{[3-(1H-tetrazol-1-yl)phenyl]oxy} acetamide, to a cell or a plurality of cells.